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=>
Uploading 09399083x.str

L1      STRUCTURE UPLOADED

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SAMPLE SEARCH INITIATED 11:32:53 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED -    131 TO ITERATE

100.0% PROCESSED     131 ITERATIONS                                  4 ANSWERS
SEARCH TIME: 00.00.03

FULL FILE PROJECTIONS:  ONLINE  **COMPLETE**
                        BATCH   **COMPLETE**
PROJECTED ITERATIONS:         1934 TO     3306
PROJECTED ANSWERS:               4 TO      200

L2            4 SEA SSS SAM L1

=> d l1
L1 HAS NO ANSWERS
L1              STR

* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY -  AVAILABLE VIA OFFLINE PRINT *

Structure attributes must be viewed using STN Express query preparation.


=> s l1 sss ful
FULL SEARCH INITIATED 11:33:50 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED -   2645 TO ITERATE

100.0% PROCESSED    2645 ITERATIONS                                 61 ANSWERS
SEARCH TIME: 00.00.03

L3           61 SEA SSS FUL L1

=> s l3
L4            3 L3

=> d l4 1-3 bib,ab,hitstr
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L4 ANSWER 1 OF 3 CAPLUS COPYRIGHT 2002 ACS
AN 2001:730744 CAPLUS
DN 135:288790
TI Pyrrolopyrimidines as tyrosine kinase inhibitors
IN Hirst, Gavin C.; Calderwood, David; Munschauer, Rainer; Arnold, Lee D.; Johnston, David N.; Rafferty, Paul
PA Basf Aktiengesellschaft, Germany
SO PCT Int. Appl., 453 pp.
CODEN: PIXXD2
DT Patent
LA English
FAN.CNT 1

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI WO 2001072751	A1	20011004	WO 2000-US8593	20000329

W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, CA, CH, CN, CR, CU, CZ, DE, DK, DM, DZ, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG, US, UZ, VN, YU, ZA, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM
RW: GH, GM, KE, LS, MW, SD, SL, SZ, TZ, UG, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG

OS MARPAT 135:288790
AB Chem. compds. having structural formula I and physiol. acceptable salts and metabolites thereof, are inhibitors of serine/threonine and tyrosine kinase activity. Several of the kinases, whose activity is inhibited by these chem. compds., are involved in immunol., hyperproliferative, or angiogenic processes. Thus, these chem. compds. can ameliorate disease states where angiogenesis or endothelial cell hyperproliferation is a factor. These compds. can be used to treat cancer and hyperproliferative disorders, rheumatoid arthritis, disorders of the immune system, transplant rejections and inflammatory disorders. All exemplified compds. significantly inhibited either FGFR, PDGFR, KDR, Tie-2, Lck, Fyn, Blk, Lyn, or Src at .ltoreq.50 .mu.M, and some significantly inhibited cdc2 at .ltoreq.50 .mu.M. In I, ring A is a six membered arom. ring or a five or six membered heteroarom. ring which is optionally substituted. L is -O-, -S-, -S(O)-, -S(O)2-, -N(R)-, -N[C(O)OR]-, -N[C(O)R]-, -N(SO2R)-, -CH2O-; -CH2S-, -CH2N(R)-, -C(NR)-;-CH2N[C(O)R]-, -CH2N[C(O)OR]-, -CH2N(SO2R)-, -CH(NHR)-, -CH[NHC(O)R]-, -CH(NHSO2R)-, -CH[NHC(O)OR]-, -CH[OC(O)R]-, -CH[OC(O)NHR]-, -CH:CH-; -C(:NOR)-, -C(O)-, -CH(OR)-, -C(O)N(R)-, -N(R)C(O)-, -N(R)S(O)-, -N(R)S(O)2-, -OC(O)N(R)-, -N(R)C(O)N(R)-, -NRC(O)O-, -S(O)N(R)-, -S(O)2N(R)-, -N[C(O)R]S(O)-, -N[C(O)R]S(O)2-, -N(R)S(O)N(R)-, -N(R)S(O)2N(R)-, -C(O)N(R)C(O)-, -S(O)N(R)C(O)-, -S(O)2N(R)C(O)-, -OS(O)N(R)-, -OS(O)2N(R)-, -N(R)S(O)O-, -N(R)S(O)2O-, -N(R)S(O)C(O)-, -N(R)S(O)2C(O)-, -SON[C(O)R]-, -SO2N[C(O)R]-, -N(R)SON(R)-, -N(R)SO2N(R)-, -C(O)O-, -N(R)P(OR')O-, -N(R)P(OR')-, -N(R)P(O)(OR')O-, -N(R)P(O)(OR')-, -N[C(O)R]P(OR')O-, -N[C(O)R]P(OR')-, -N[C(O)R]P(O)(OR')O-, -N[C(O)R]P(OR')-, -CH(R)S(O)-, or -CH(R)S(O)2-. L is also -CH(R)N[C(O)OR]-, -CH(R)N[C(O)R]-, -CH(R)N(SO2R), -CH(R)O-, -CH(R)S-, -CH(R)N(R)-, -CH(R)N[C(O)R]-, - CH(R)N[C(O)OR]-, -CH(R)N(SO2R)-, -CH(R)C(:NOR)-, -CH(R)C(O)-, -CH(R)CH(OR)-, -CH(R)C(O)N(R)-, -CH(R)N(R)C(O)-, -CH(R)N(R)S(O)-, -CH(R)N(R)S(O)2-, -CH(R)OC(O)N(R)-, -CH(R)N(R)C(O)N(R)-, -CH(R)N(R)C(O)O-, -CH(R)S(O)N(R)-, -CH(R)S(O)2N(R)-, -CH(R)N[C(O)R]S(O)-, -CH(R)N[C(O)R]S(O)2-, -CH(R)N(R)S(O)N(R)-, -CH(R)N(R)S(O)2N(R)-, -CH(R)C(O)N(R)C(O)-, -CH(R)S(O)N(R)C(O)-, -CH(R)S(O)2N(R)C(O)-, -CH(R)OS(O)N(R)-, -CH(R)OS(O)2N(R)-,

-CH(R)N(R)S(O)O-, -CH(R)N(R)S(O)2O-, -CH(R)N(R)S(O)C(O)-, -CH(R)N(R)S(O)2C(O)-, -CH(R)SON[C(O)R]-, -CH(R)S(O)2N[C(O)R]-, -CH(R)N(R)SON(R)-, -CH(R)N(R)S(O)2N(R)-; -CH(R)C(O)O-, -CH(R)N(R)P(OR')O-, -CH(R)N(R)P(OR')-, -CH(R)N(R)P(O)(OR')O-, -CH(R)N(R)P(O)(OR')-, -CH(R)N[C(O)R]P(OR')O-, -CH(R)N[C(O)R]P(OR')-, -CH(R)N[C(O)R]P(O)(OR')O- or -CH(R)N[C(O)R]P(OR')-. In L, each R and R' is, independently, -H, acyl, substituted or unsubstituted aliph., arom., arylalkyl, heteroarom., cycloalkyl or arylalkyl; or L is -RbN(R)S(O)2-, -RbN(R)P(O)-, or -RbN(R)P(O)O-, wherein Rb is an alkylene group which when taken together with the sulfonamide, phosphinamide, or phosphonamide group to which it is bound forms a five or six membered ring fused to ring A; or L is II (X = O or nil; Y = O or nil) or III (Y = O, nil) wherein R85 taken together with the phosphinamide, or phosphonamide is a 5-, 6-, or 7-membered, arom., heteroarom. or heterocycloalkyl ring system. G is a direct bond, -(CH2)j- (j = 1-6), C2-C6-alkenylene, C3-C8-cycloalkylene or C1-C6-oxaalkylene group. R1 is substituted or optionally substituted aliph., cycloalkyl, bicycloalkyl, cycloalkenyl, arom., heteroarom., heteroaralkyl, heterocycloalkyl, heterobicycloalkyl, alkylamido, arylamido, -S(O)2-alkyl, -S(O)2-cycloalkyl, -C(O)alkyl, or -B-E, wherein B is substituted or unsubstituted cycloalkyl, heterocycloalkyl, arom., heteroarom., alkylene, aminoalkyl, alkylenecarbonyl, or aminoalkylcarbonyl and E is substituted or unsubstituted azacycloalkyl, azacycloalkylcarbonyl, azacycloalkylsulfonyl, azacycloalkylalkyl, heteroaryl, heteroarylcarbonyl, heteroarylsulfonyl, heteroaralkyl, alkyl sulfonamido, aryl sulfonamido, bicycloalkyl, ureido, thioureido or aryl. R2 is -H or substituted or unsubstituted aliph., cycloalkyl, halogen, -OH, cyano, arom., heteroarom., heterocycloalkyl, aralkyl, heteroaralkyl, -(CH2)0-3NR4R5, or -(CH2)0-3C(O)NR4R5. R3 is substituted or unsubstituted aliph., alkenyl, cycloalkyl, arom., heteroarom., or heterocycloalkyl with provisos. R4, R5 and the N atom together form a 3, 4, 5, 6 or 7-membered, substituted or unsubstituted heterocycloalkyl, heterobicycloalkyl or heteroarom.; or R4 and R5 are each, independently, -H, azabicycloalkyl, heterocycloalkyl, substituted or unsubstituted alkyl or Y-Z; Y is -C(O)-, -(CH2)p-, -S(O)2-, -C(O)O-, -SO2NH-, -CONH-, -(CH2)pO-, -(CH2)pNH-, -(CH2)pS-, -(CH2)pS(O)-, and -(CH2)pS(O)2-; p = 0-6; and Z is -H, or substituted or unsubstituted alkyl, amino, aryl, heteroaryl or heterocycloalkyl. 546 Example prepns. are included. For example, addn. of piperidine to 4-[4-amino-5-(4-phenoxyphenyl)-7H-pyrrolo[2,3-d]pyrimidin-7-yl]cyclohexanone in DCE and AcOH, followed by treatment with Na[(AcO)3BH], workup and chromatog., gave cis- and trans-IV.

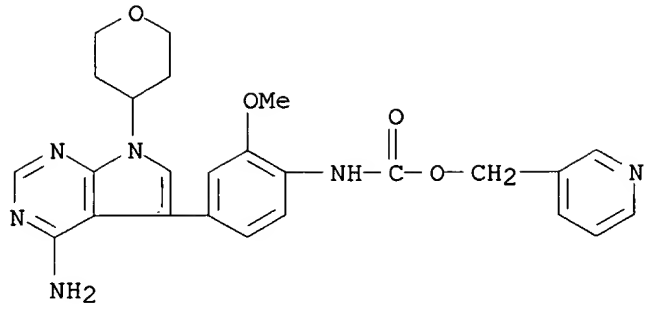
IT **262444-13-3P 262444-14-4P 262444-16-6P 262444-17-7P**

RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); RCT (Reactant); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); RACT (Reactant or reagent); USES (Uses)

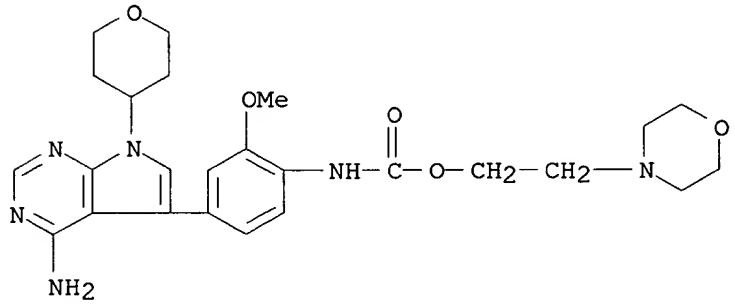
(intermediate; prepn. of pyrrolopyrimidinamines as protein kinase inhibitors)

RN 262444-13-3 CAPLUS

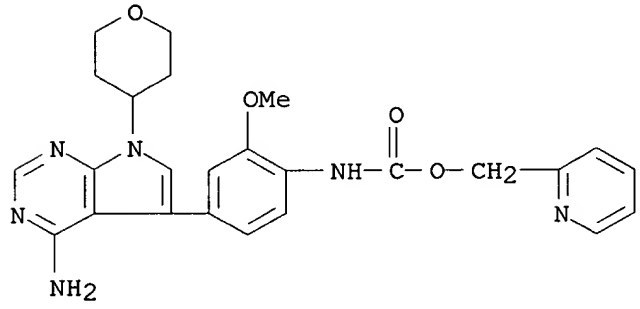
CN Carbamic acid, [4-[4-amino-7-(tetrahydro-2H-pyran-4-yl)-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-, 3-pyridinylmethyl ester (9CI) (CA INDEX NAME)



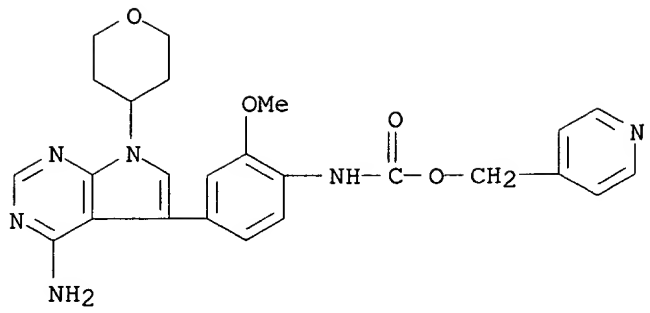
RN 262444-14-4 CAPLUS
CN Carbamic acid, [4-[4-amino-7-(tetrahydro-2H-pyran-4-yl)-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-, 2-(4-morpholinyl)ethyl ester (9CI) (CA INDEX NAME)



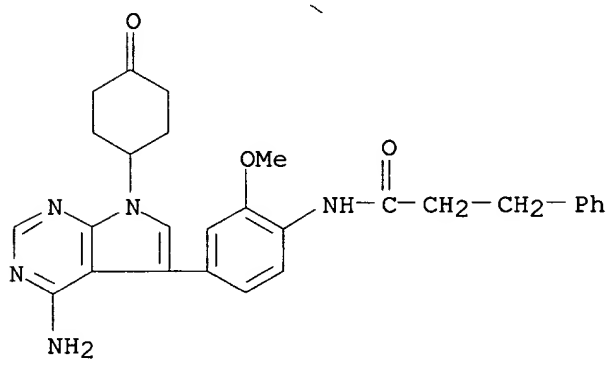
RN 262444-16-6 CAPLUS
CN Carbamic acid, [4-[4-amino-7-(tetrahydro-2H-pyran-4-yl)-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-, 2-pyridinylmethyl ester (9CI) (CA INDEX NAME)



RN 262444-17-7 CAPLUS
CN Carbamic acid, [4-[4-amino-7-(tetrahydro-2H-pyran-4-yl)-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-, 4-pyridinylmethyl ester (9CI) (CA INDEX NAME)



IT **262444-51-9P**
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)
(intermediate; prepn. of pyrrolopyrimidinamines as protein kinase inhibitors)
RN 262444-51-9 CAPLUS
CN Benzenepropanamide, N-[4-[4-amino-7-(4-oxocyclohexyl)-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]- (9CI) (CA INDEX NAME)

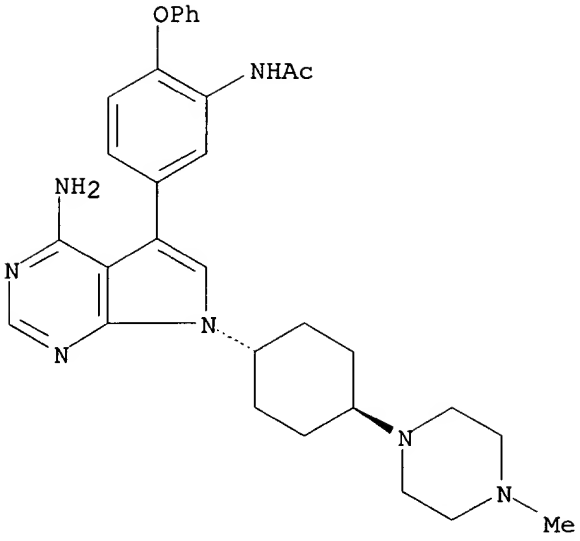


IT **364354-19-8P 364354-56-3P 364354-57-4P**
364354-68-7P 364356-13-8P
RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
(prepn. of pyrrolopyrimidinamines as protein kinase inhibitors)
RN 364354-19-8 CAPLUS
CN Acetamide, N-[5-[4-amino-7-[trans-4-(4-methyl-1-piperazinyl)cyclohexyl]-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-phenoxyphenyl]-, monoacetate (9CI) (CA INDEX NAME)

CM 1

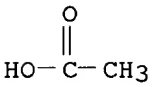
CRN 364354-18-7
CMF C31 H37 N7 O2

Relative stereochemistry.



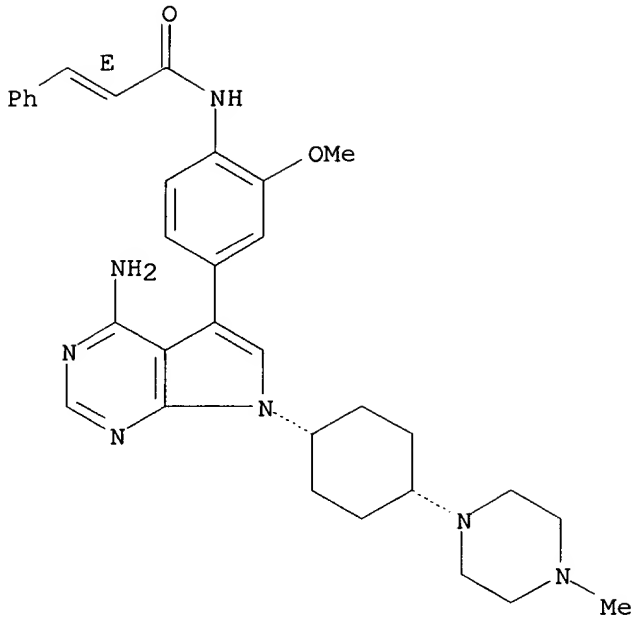
CM 2

CRN 64-19-7
CMF C2 H4 O2



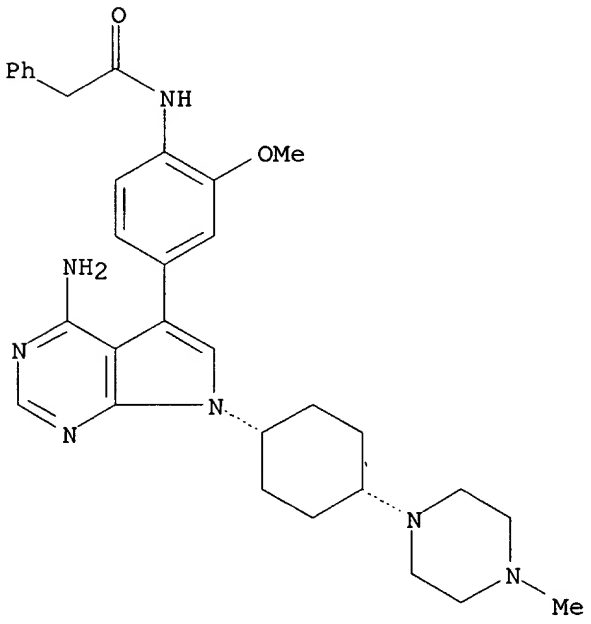
RN 364354-56-3 CAPLUS
CN 2-Propenamide, N-[4-[4-amino-7-[cis-4-(4-methyl-1-piperazinyl)cyclohexyl]-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-3-phenyl-, (2E)- (9CI) (CA INDEX NAME)

Relative stereochemistry.
Double bond geometry as shown.



RN 364354-57-4 CAPLUS
CN Benzeneacetamide, N-[4-[4-amino-7-[cis-4-(4-methyl-1-piperazinyl)cyclohexyl]-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-(9CI) (CA INDEX NAME)

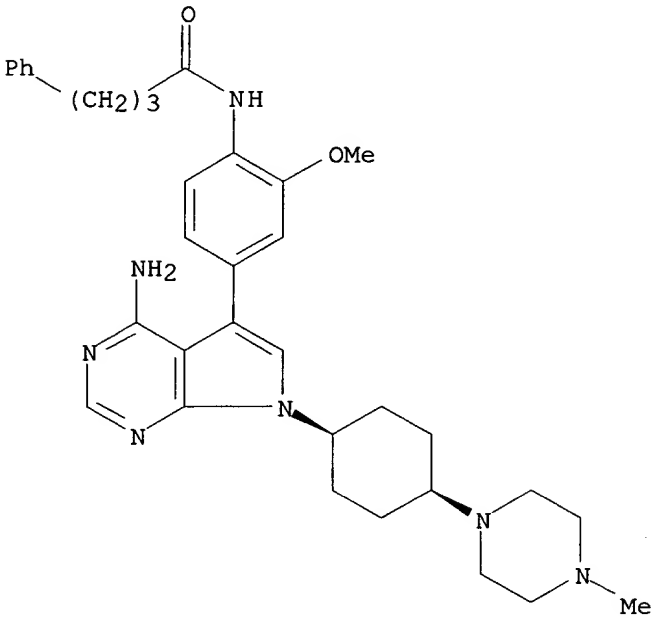
Relative stereochemistry.



RN 364354-68-7 CAPLUS
CN Benzenebutanamide, N-[4-[4-amino-7-[cis-4-(4-methyl-1-

piperazinyl)cyclohexyl]-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-
(9CI) (CA INDEX NAME)

Relative stereochemistry.



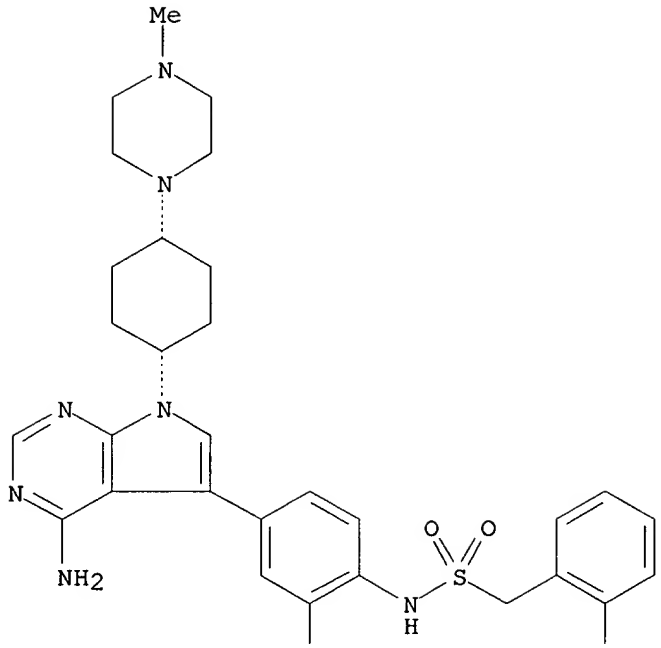
RN 364356-13-8 CAPLUS
CN Benzenemethanesulfonamide, N-[4-[4-amino-7-[cis-4-(4-methyl-1-piperazinyl)cyclohexyl]-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-fluorophenyl]-2-nitro-, (2Z)-2-butenedioate (1:2) (9CI) (CA INDEX NAME)

CM 1

CRN 262443-68-5
CMF C30 H35 F N8 O4 S

Relative stereochemistry.

PAGE 1-A

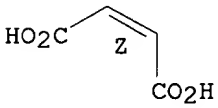


PAGE 2-A

CM 2

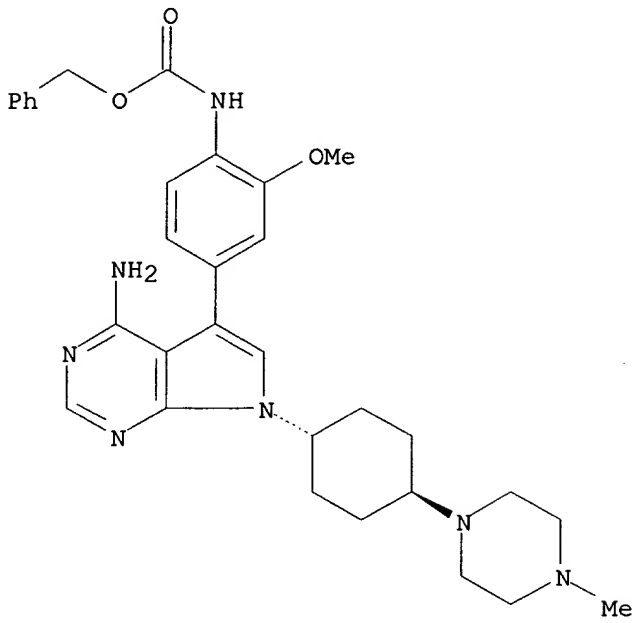
CRN 110-16-7
CMF C4 H4 O4
CDES 2:Z

Double bond geometry as shown.



IT **262442-79-5**
RL: RCT (Reactant); RACT (Reactant or reagent)
(reactant; prepn. of pyrrolopyrimidinamines as protein kinase inhibitors)
RN 262442-79-5 CAPLUS
CN Carbamic acid, [4-[4-amino-7-[trans-4-(4-methyl-1-piperazinyl)cyclohexyl]-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-, phenylmethyl ester (9CI) (CA INDEX NAME)

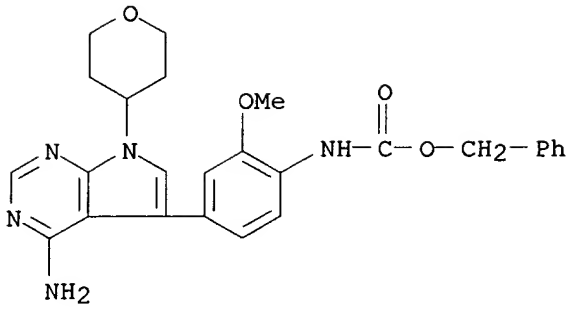
Relative stereochemistry.



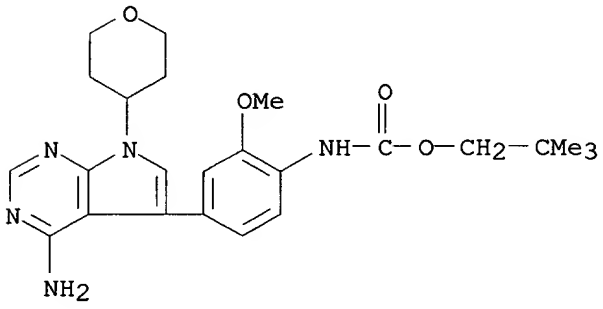
IT **262439-86-1P 262439-87-2P 262439-90-7P**
262439-91-8P 262439-92-9P 262439-96-3P
262439-97-4P 262439-98-5P 262439-99-6P
262440-00-6P 262440-04-0P 262440-06-2P
262442-22-8P 262442-32-0P 262442-33-1P
262442-39-7P 262442-47-7P 262442-80-8P
262443-69-6P 262443-91-4P 262444-01-9P
262444-02-0P 262444-03-1P 262444-68-8P
262445-41-0P
RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
(target compd.; prepn. of pyrrolopyrimidinamines as protein kinase inhibitors)

RN 262439-86-1 CAPLUS

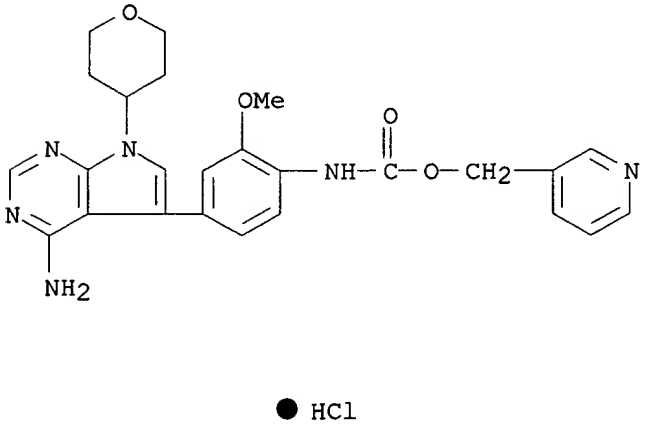
CN Carbamic acid, [4-[4-amino-7-(tetrahydro-2H-pyran-4-yl)-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-, phenylmethyl ester (9CI) (CA INDEX NAME)



RN 262439-87-2 CAPLUS
CN Carbamic acid, [4-[4-amino-7-(tetrahydro-2H-pyran-4-yl)-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-, 2,2-dimethylpropyl ester (9CI) (CA INDEX NAME)



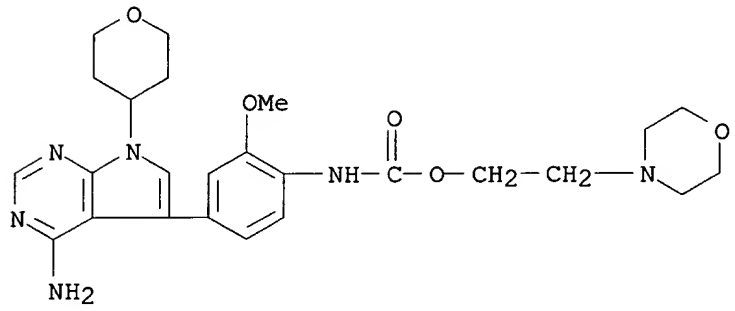
RN 262439-90-7 CAPLUS
CN Carbamic acid, [4-[4-amino-7-(tetrahydro-2H-pyran-4-yl)-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-, 3-pyridinylmethyl ester, monohydrochloride (9CI) (CA INDEX NAME)



● HCl

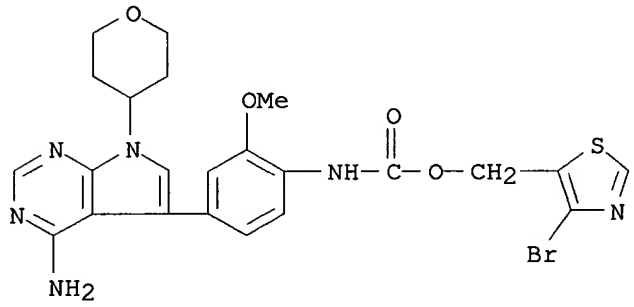
RN 262439-91-8 CAPLUS
CN Carbamic acid, [4-[4-amino-7-(tetrahydro-2H-pyran-4-yl)-7H-pyrrolo[2,3-

d]pyrimidin-5-yl]-2-methoxyphenyl]-, 2-(4-morpholinyl)ethyl ester, monohydrochloride (9CI) (CA INDEX NAME)

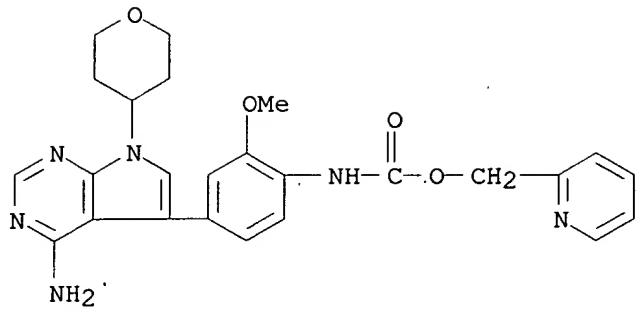


● HCl

RN 262439-92-9 CAPLUS
CN Carbamic acid, [4-[4-amino-7-(tetrahydro-2H-pyran-4-yl)-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-, (4-bromo-5-thiazolyl)methyl ester (9CI) (CA INDEX NAME)

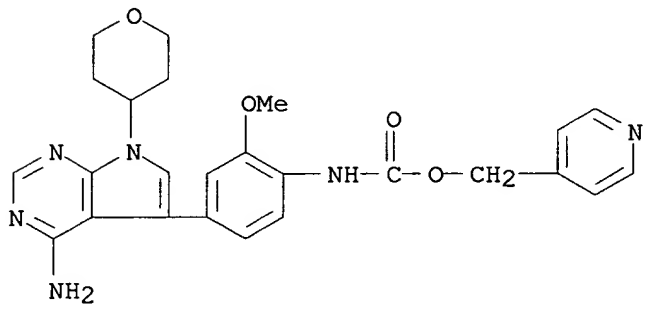


RN 262439-96-3 CAPLUS
CN Carbamic acid, [4-[4-amino-7-(tetrahydro-2H-pyran-4-yl)-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-, 2-pyridinylmethyl ester, monohydrochloride (9CI) (CA INDEX NAME)



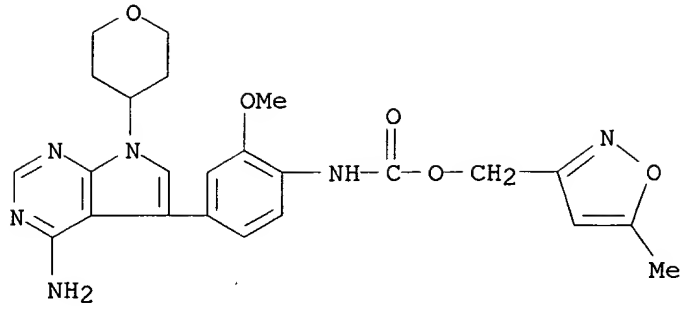
● HCl

RN 262439-97-4 CAPLUS
CN Carbamic acid, [4-[4-amino-7-(tetrahydro-2H-pyran-4-yl)-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-, 4-pyridinylmethyl ester, monohydrochloride (9CI) (CA INDEX NAME)



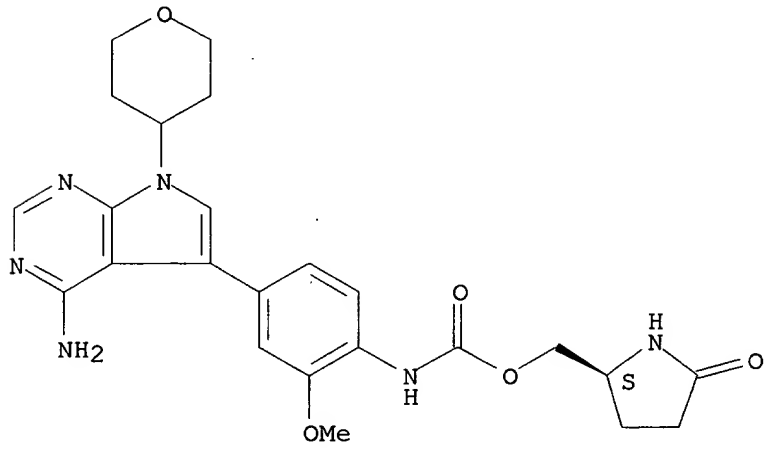
● HCl

RN 262439-98-5 CAPLUS
CN Carbamic acid, [4-[4-amino-7-(tetrahydro-2H-pyran-4-yl)-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-, (5-methyl-3-isoxazolyl)methyl ester (9CI) (CA INDEX NAME)

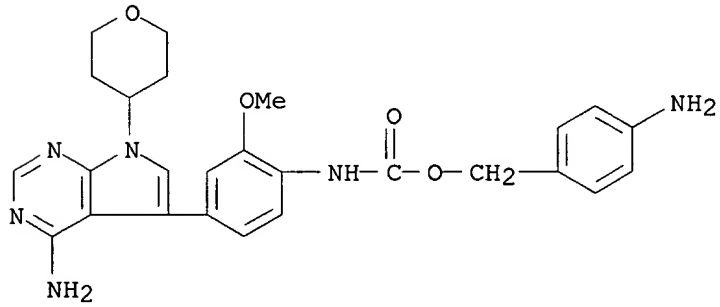


RN 262439-99-6 CAPLUS
CN Carbamic acid, [4-[4-amino-7-(tetrahydro-2H-pyran-4-yl)-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-, [(2S)-5-oxo-2-pyrrolidinyl]methyl ester (9CI) (CA INDEX NAME)

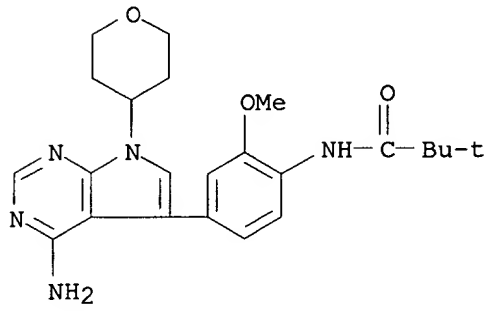
Absolute stereochemistry.



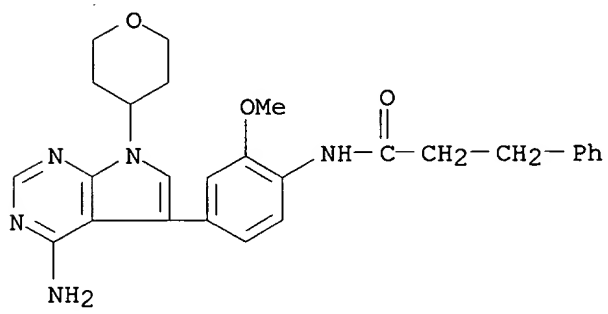
RN 262440-00-6 CAPLUS
CN Carbamic acid, [4-[4-amino-7-(tetrahydro-2H-pyran-4-yl)-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-, (4-aminophenyl)methyl ester (9CI) (CA INDEX NAME)



RN 262440-04-0 CAPLUS
CN Propanamide, N-[4-[4-amino-7-(tetrahydro-2H-pyran-4-yl)-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-2,2-dimethyl- (9CI) (CA INDEX NAME)

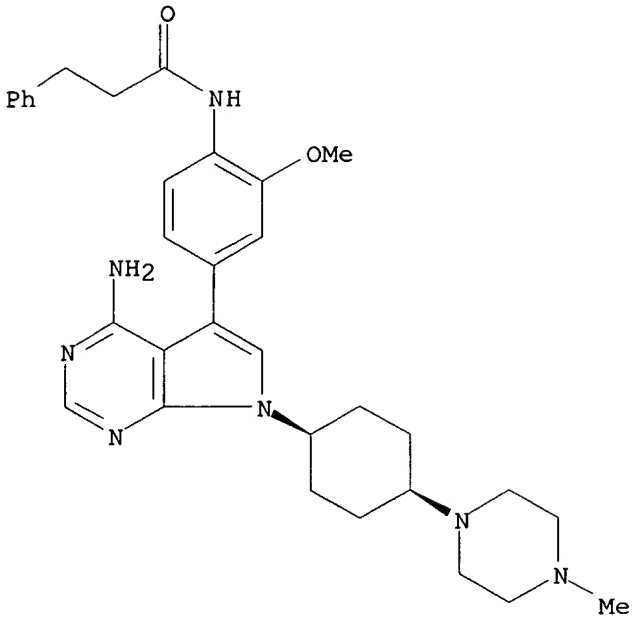


RN 262440-06-2 CAPLUS
CN Benzenepropanamide, N-[4-[4-amino-7-(tetrahydro-2H-pyran-4-yl)-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]- (9CI) (CA INDEX NAME)



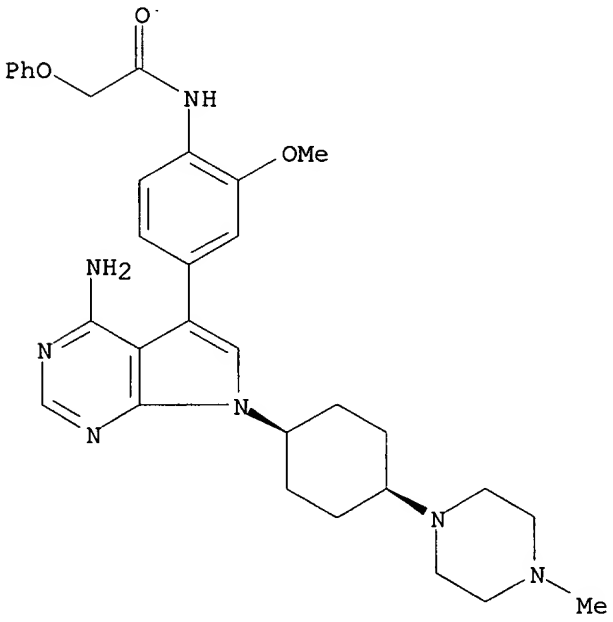
RN 262442-22-8 CAPLUS
CN Benzenepropanamide, N-[4-[4-amino-7-[cis-4-(4-methyl-1-piperazinyl)cyclohexyl]-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]- (9CI) (CA INDEX NAME)

Relative stereochemistry.



RN 262442-32-0 CAPLUS
CN Acetamide, N-[4-[4-amino-7-[cis-4-(4-methyl-1-piperazinyl)cyclohexyl]-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-2-phenoxy- (9CI) (CA INDEX NAME)

Relative stereochemistry.



RN 262442-33-1 CAPLUS
CN Acetamide, N-[4-[4-amino-7-[cis-4-(4-methyl-1-piperazinyl)cyclohexyl]-7H-

pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-2-(4-chlorophenoxy)- (9CI) (CA INDEX NAME)

Relative stereochemistry.

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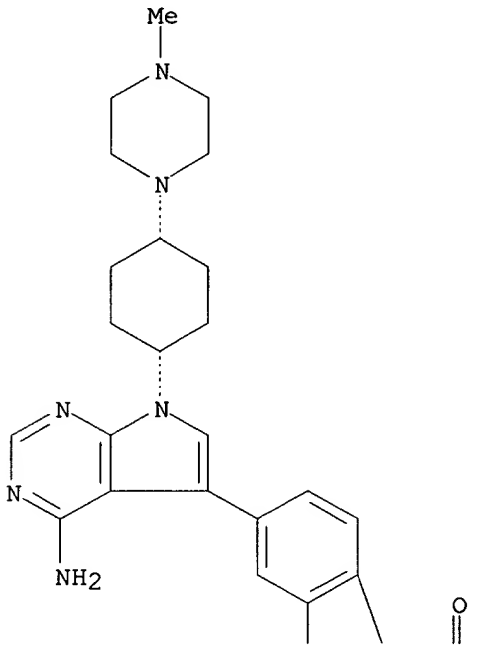
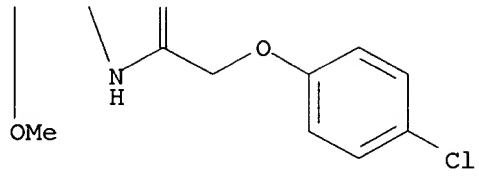


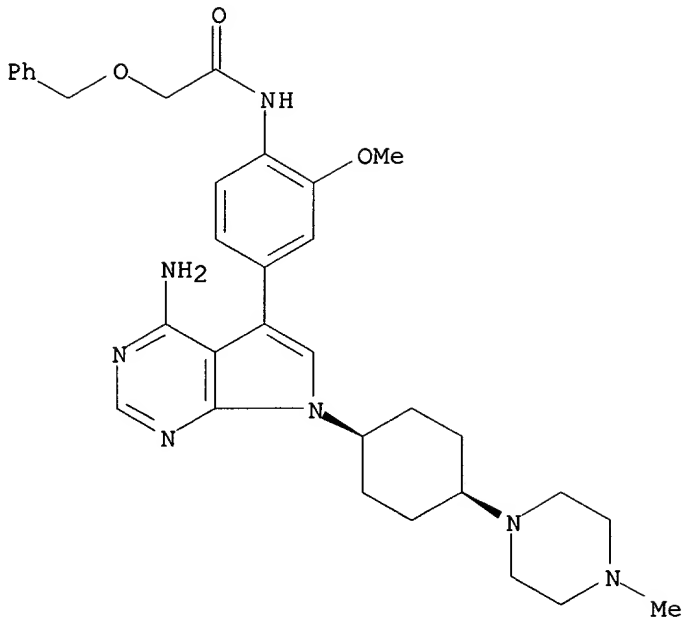
PAGE 2-A



RN 262442-39-7 CAPLUS

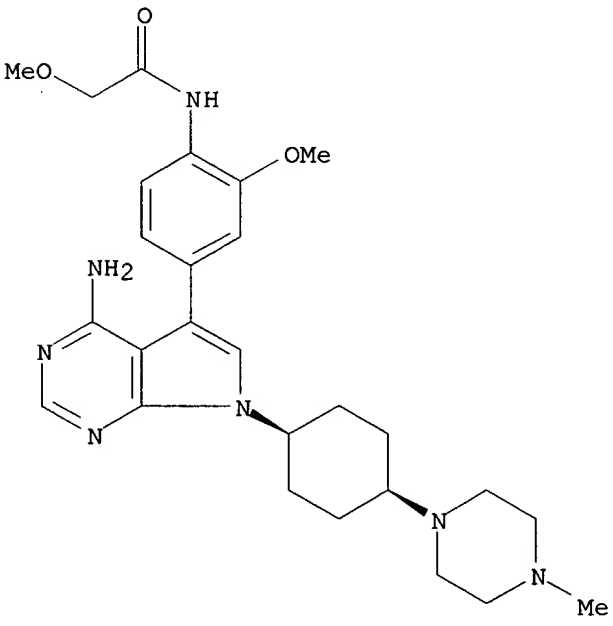
CN Acetamide, N-[4-[4-amino-7-[cis-4-(4-methyl-1-piperazinyl)cyclohexyl]-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-2-(phenylmethoxy)- (9CI) (CA INDEX NAME)

Relative stereochemistry.



RN 262442-47-7 CAPLUS
CN Acetamide, N-[4-[4-amino-7-[cis-4-(4-methyl-1-piperazinyl)cyclohexyl]-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-2-methoxy- (9CI) (CA INDEX NAME)

Relative stereochemistry.



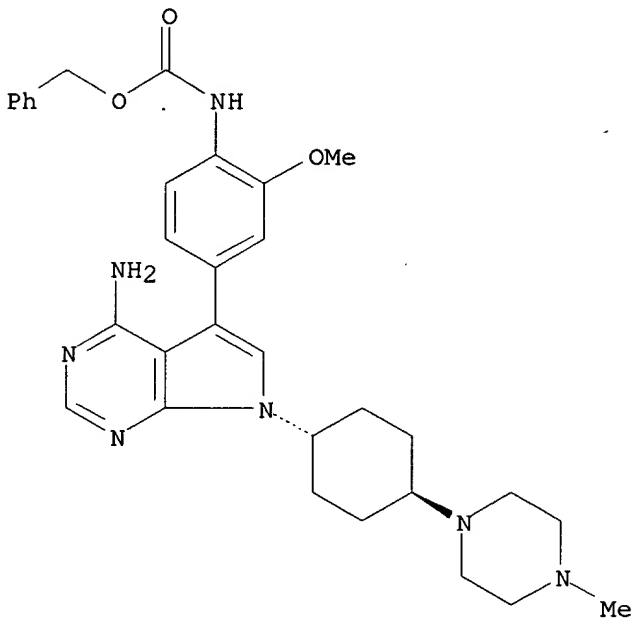
RN 262442-80-8 CAPLUS
CN Carbamic acid, [4-[4-amino-7-[trans-4-(4-methyl-1-piperazinyl)cyclohexyl]-

7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-, phenylmethyl ester, (2Z)-2-butenedioate (1:3) (9CI) (CA INDEX NAME)

CM 1

CRN 262442-79-5
CMF C32 H39 N7 O3

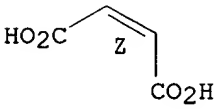
Relative stereochemistry.



CM 2

CRN 110-16-7
CMF C4 H4 O4
CDES 2:Z

Double bond geometry as shown.



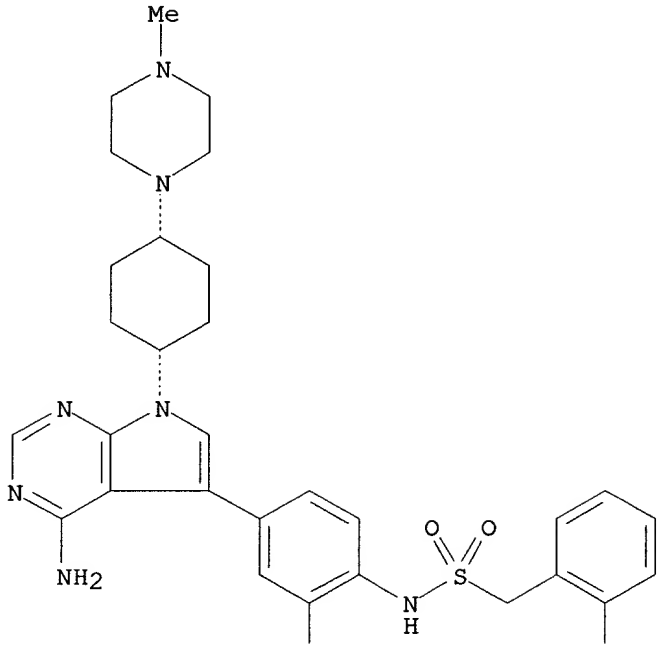
RN 262443-69-6 CAPLUS
CN Benzenemethanesulfonamide, N-[4-[4-amino-7-[cis-4-(4-methyl-1-piperazinyl)cyclohexyl]-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-fluorophenyl]-2-nitro-, (2Z)-2-butenedioate (1:3) (9CI) (CA INDEX NAME)

CM 1

CRN 262443-68-5
CMF C30 H35 F N8 O4 S

Relative stereochemistry.

PAGE 1-A



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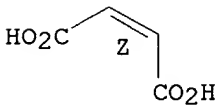
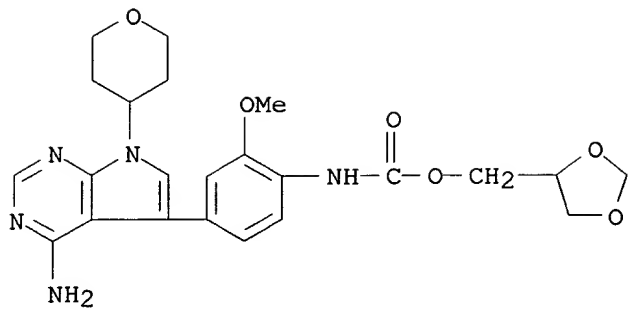
CM 2

CRN 110-16-7
CMF C4 H4 O4
CDES 2:Z

Double bond geometry as shown.



RN 262443-91-4 CAPLUS
CN Carbamic acid, [4-[4-amino-7-(tetrahydro-2H-pyran-4-yl)-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-, 1,3-dioxolan-4-ylmethyl ester (9CI) (CA INDEX NAME)

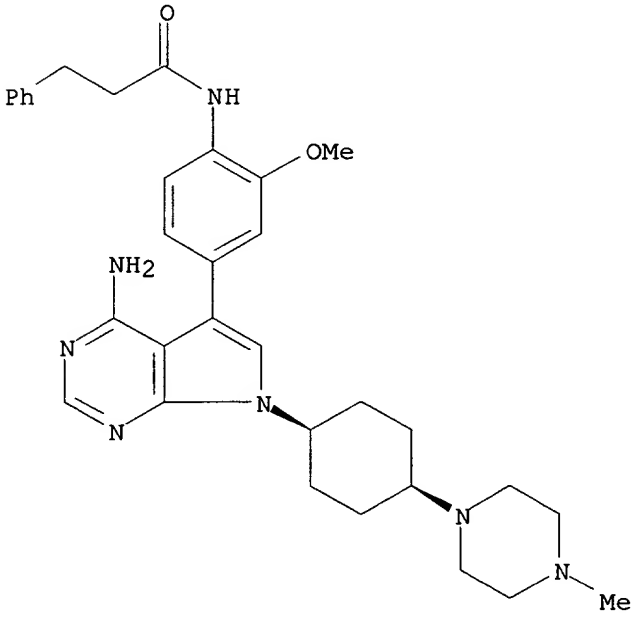


RN 262444-01-9 CAPLUS
CN Benzenepropanamide, N-[4-[4-amino-7-[cis-4-(4-methyl-1-piperazinyl)cyclohexyl]-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-, (2Z)-2-butenedioate (1:3) (9CI) (CA INDEX NAME)

CM 1

CRN 262442-22-8
CMF C33 H41 N7 O2

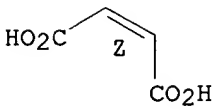
Relative stereochemistry.



CM 2

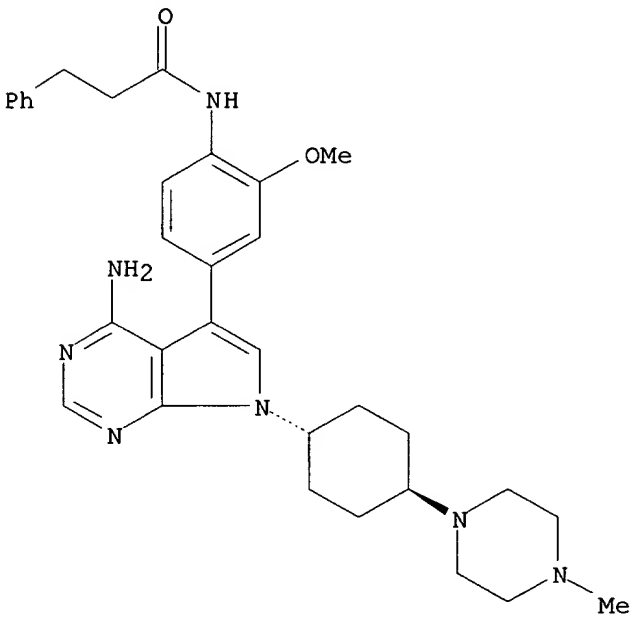
CRN 110-16-7
CMF C4 H4 O4
CDES 2:Z

Double bond geometry as shown.



RN 262444-02-0 CAPLUS
CN Benzenepropanamide, N-[4-[4-amino-7-[trans-4-(4-methyl-1-piperazinyl)cyclohexyl]-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-(9CI) (CA INDEX NAME)

Relative stereochemistry.

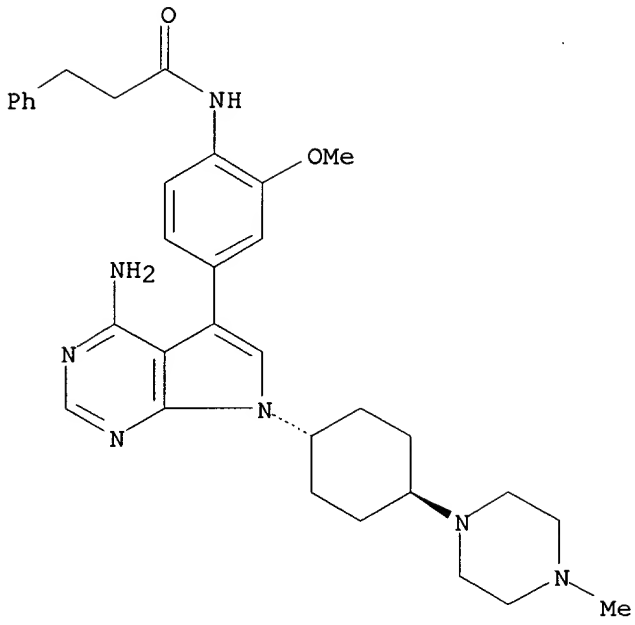


RN 262444-03-1 CAPLUS
CN Benzenepropanamide, N-[4-[4-amino-7-[trans-4-(4-methyl-1-piperazinyl)cyclohexyl]-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-, (2Z)-2-butenedioate (1:3) (9CI) (CA INDEX NAME)

CM 1

CRN 262444-02-0
CMF C33 H41 N7 O2

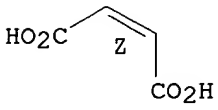
Relative stereochemistry.



CM 2

CRN 110-16-7
CMF C4 H4 O4
CDES 2:Z

Double bond geometry as shown.

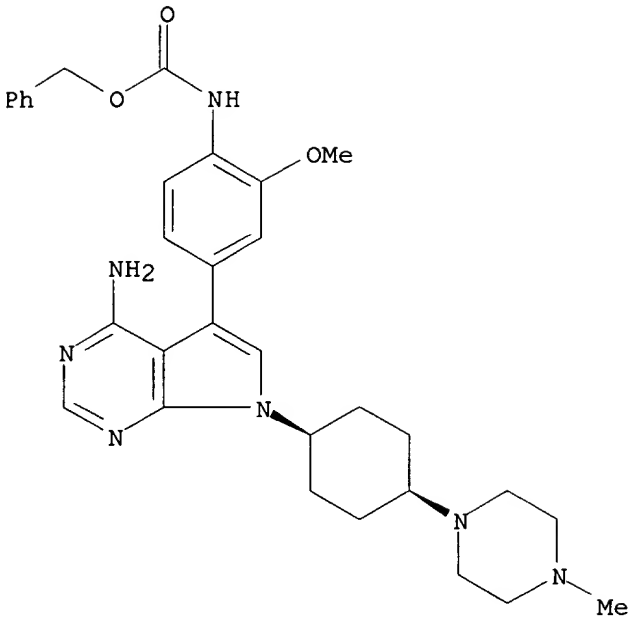


RN 262444-68-8 CAPLUS
CN Carbamic acid, [4-[4-amino-7-[cis-4-(4-methyl-1-piperazinyl)cyclohexyl]-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-, phenylmethyl ester, (2Z)-2-butenedioate (1:3) (9CI) (CA INDEX NAME)

CM 1

CRN 262444-67-7
CMF C32 H39 N7 O3

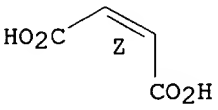
Relative stereochemistry.



CM 2

CRN 110-16-7
CMF C4 H4 O4
CDES 2:Z

Double bond geometry as shown.



RN 262445-41-0 CAPLUS
CN Benzenemethanesulfonamide, N-[4-[4-amino-7-[trans-4-(4-methyl-1-piperazinyl)cyclohexyl]-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-fluorophenyl]-2-nitro-, (2Z)-2-butenedioate (1:3) (9CI) (CA INDEX NAME)

CM 1

CRN 262445-40-9
CMF C30 H35 F N8 O4 S

Relative stereochemistry.

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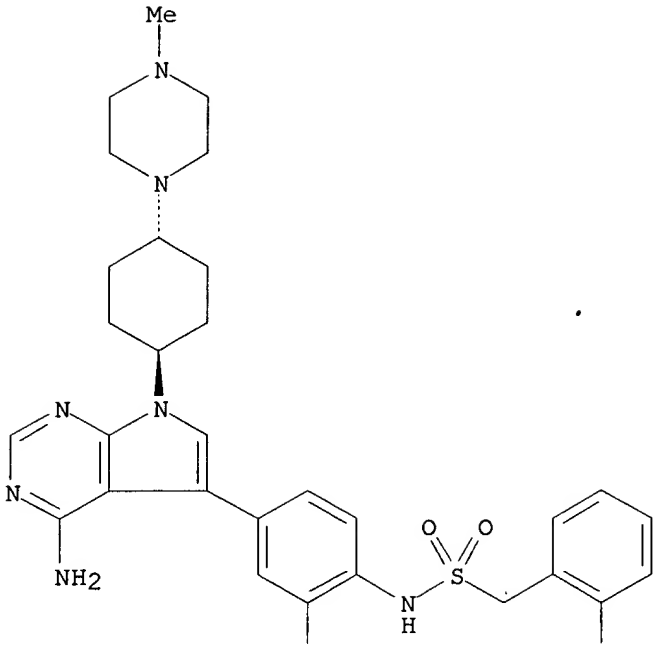

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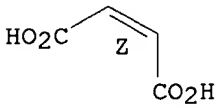
CM 2

CRN 110-16-7
CMF C4 H4 O4
CDES 2:Z

Double bond geometry as shown.



RE.CNT 14 THERE ARE 14 CITED REFERENCES AVAILABLE FOR THIS RECORD
ALL CITATIONS AVAILABLE IN THE RE FORMAT

L4 ANSWER 2 OF 3 CAPLUS COPYRIGHT 2002 ACS
AN 2000:210172 CAPLUS
DN 132:251160
TI Preparation of pyrrolopyrimidines as protein kinase inhibitors
IN Hirst, Gavin C.; Calderwood, David; Wishart, Neil; Ritter, Kurt; Arnold, Lee D.
PA Basf A.-G., Germany
SO PCT Int. Appl., 304 pp.
CODEN: PIXXD2
DT Patent
LA English
FAN.CNT 1

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI WO 2000017203	A1	20000330	WO 1999-US21560	19990917
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RW: GH, GM, KE, LS, MW, SD, SL, SZ, TZ, UG, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG				
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EP 1114053	A1	20010711	EP 1999-969415	19990917
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BR 9913887	A	20011023	BR 1999-13887	19990917
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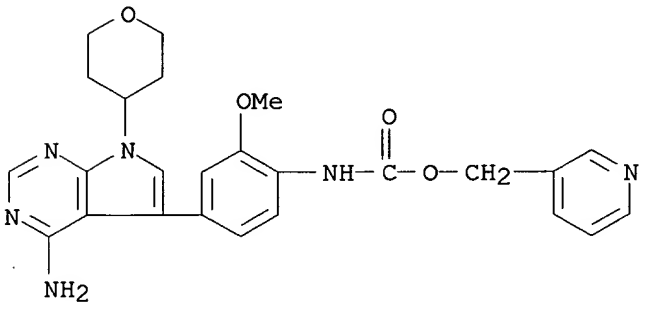
PRAI US 1998-100832P	P	19980918		
US 1998-100833P	P	19980918		
US 1998-100834P	P	19980918		
US 1998-100946P	P	19980918		
WO 1999-US21560	W	19990917		

OS MARPAT 132:251160
AB 7H-Pyrrolo[2,3-d]pyrimidin-4-amines (I) [wherein A = (un)substituted 6-membered arom. ring or 5- or 6-membered heteroarom. ring; L = RbN(R)S(O)2, RbN(R)P(O), or RbN(R)P(O)O, where Rb = alkylene group which when taken together with the sulfonamide, phosphinamide or phosphonamide group to which it is bound forms a 5- or 6-membered ring fused to ring A, or L = 5-, 6-, or 7-membered (oxa)azaphosphaarom. or (oxa)azaphosphacycloalkyl ring; R = H, acyl, or (un)substituted aliph., (hetero)arom., or cycloalkyl; R1 = (un)substituted (hetero)cyclic, (hetero)arom., amido, acyl, or (cyclo)alkylsulfonyl; R2 = H, halo, OH, CN, (un)substituted aliph., cycloalkyl, (hetero)arom., (hetero)aralkyl, amino, or amido; R3 (un)substituted aliph., alkenyl, (hetero)cycloalkyl, or (hetero)arom.; n = 0-6], and physiol. acceptable salts and metabolites thereof, were prepd. For example, addn. of piperidine to 4-[4-amino-5-(4-phenoxyphenyl)-7H-pyrrolo[2,3-d]pyrimidin-7-yl]cyclohexanone in DCE and AcOH, followed by workup and chromatog., gave cis- and trans-II. I inhibit serine/threonine and tyrosine kinase activity, which are involved in immunol., hyperproliferative, and angiogenic processes. All exemplified compds. significantly inhibited either FGFR, PDGFR, KDR, Tie-2, Lck, Fyn, Blk, Lyn, or Src at concns. of .ltoreq. 50 .mu.M, and some significantly inhibited cdc2 at concns. of 50 .ltoreq. .mu.M. Thus, these compds. are useful in the treatment of cancer and hyperproliferative disorders, rheumatoid arthritis, disorders of the

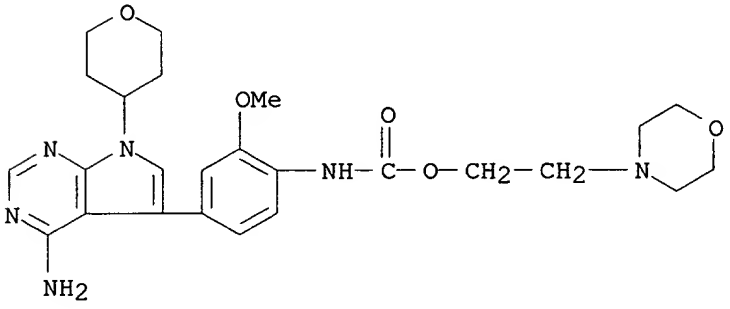
immune system, transplant rejections, and inflammatory disorders.

IT **262444-13-3P 262444-14-4P 262444-16-6P 262444-17-7P**
RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); RCT (Reactant); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); RACT (Reactant or reagent); USES (Uses)
(intermediate; prepn. of 7H-pyrrolo[2,3-d]pyrimidin-4-amines as protein kinase inhibitors)

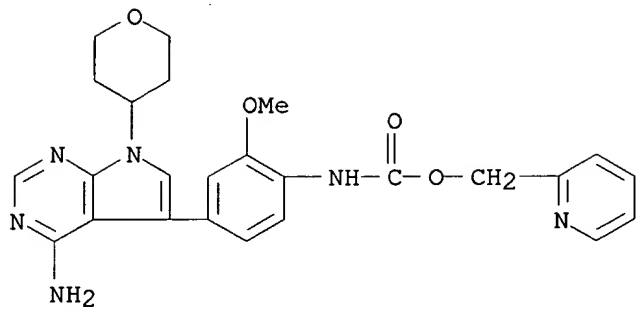
RN 262444-13-3 CAPLUS
CN Carbamic acid, [4-[4-amino-7-(tetrahydro-2H-pyran-4-yl)-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-, 3-pyridinylmethyl ester (9CI) (CA INDEX NAME)



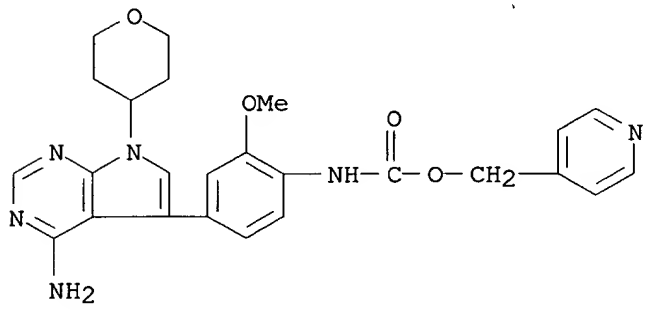
RN 262444-14-4 CAPLUS
CN Carbamic acid, [4-[4-amino-7-(tetrahydro-2H-pyran-4-yl)-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-, 2-(4-morpholinyl)ethyl ester (9CI) (CA INDEX NAME)



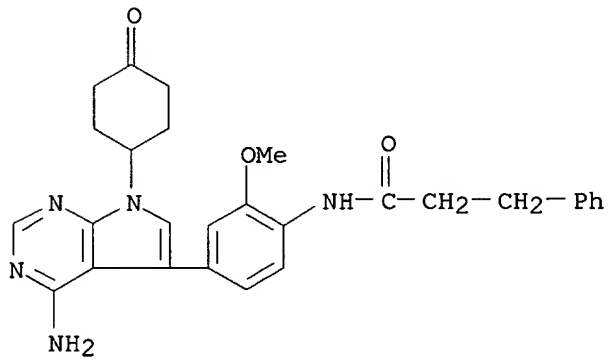
RN 262444-16-6 CAPLUS
CN Carbamic acid, [4-[4-amino-7-(tetrahydro-2H-pyran-4-yl)-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-, 2-pyridinylmethyl ester (9CI) (CA INDEX NAME)



RN 262444-17-7 CAPLUS
CN Carbamic acid, [4-[4-amino-7-(tetrahydro-2H-pyran-4-yl)-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-, 4-pyridinylmethyl ester (9CI) (CA INDEX NAME)

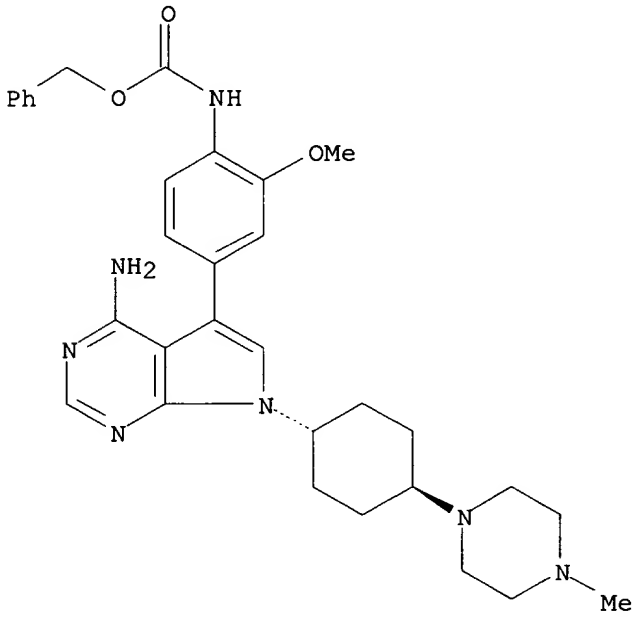


IT **262444-51-9P**
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)
(intermediate; prepn. of 7H-pyrrolo[2,3-d]pyrimidin-4-amines as protein kinase inhibitors)
RN 262444-51-9 CAPLUS
CN Benzenepropanamide, N-[4-[4-amino-7-(4-oxocyclohexyl)-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]- (9CI) (CA INDEX NAME)

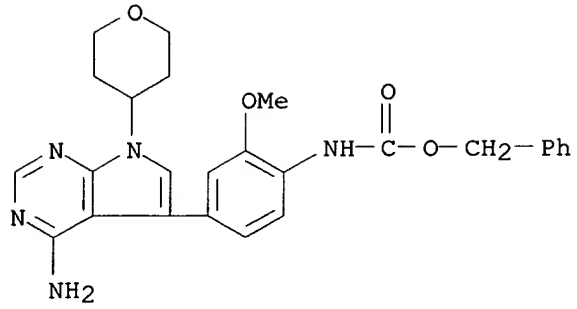


IT **262442-79-5**
RL: RCT (Reactant); RACT (Reactant or reagent)
(reactant; prepn. of 7H-pyrrolo[2,3-d]pyrimidin-4-amines as protein kinase inhibitors)
RN 262442-79-5 CAPLUS
CN Carbamic acid, [4-[4-amino-7-[trans-4-(4-methyl-1-piperazinyl)cyclohexyl]-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-, phenylmethyl ester (9CI) (CA INDEX NAME)

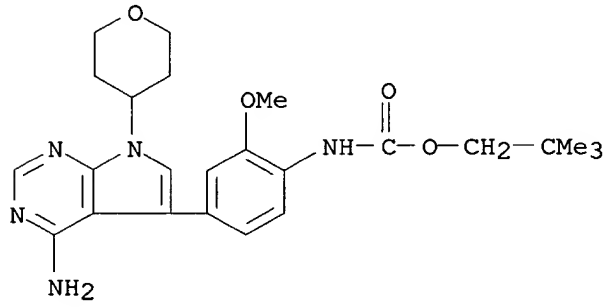
Relative stereochemistry.



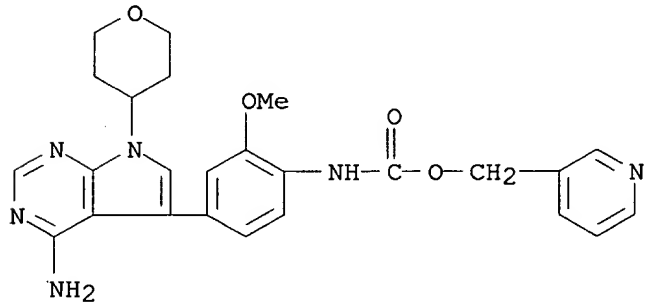
IT **262439-86-1P 262439-87-2P 262439-90-7P**
262439-91-8P 262439-92-9P 262439-96-3P
262439-97-4P 262439-98-5P 262439-99-6P
262440-00-6P 262440-04-0P 262440-06-2P
262442-22-8P 262442-32-0P 262442-33-1P
262442-39-7P 262442-47-7P 262442-80-8P
262443-69-6P 262443-91-4P 262444-01-9P
262444-02-0P 262444-03-1P 262444-68-8P
262445-41-0P
RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
(target compd.; prepn. of 7H-pyrrolo[2,3-d]pyrimidin-4-amines as protein kinase inhibitors)
RN 262439-86-1 CAPLUS
CN Carbamic acid, [4-[4-amino-7-(tetrahydro-2H-pyran-4-yl)-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-, phenylmethyl ester (9CI) (CA INDEX NAME)



RN 262439-87-2 CAPLUS
CN Carbamic acid, [4-[4-amino-7-(tetrahydro-2H-pyran-4-yl)-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-, 2,2-dimethylpropyl ester (9CI) (CA INDEX NAME)



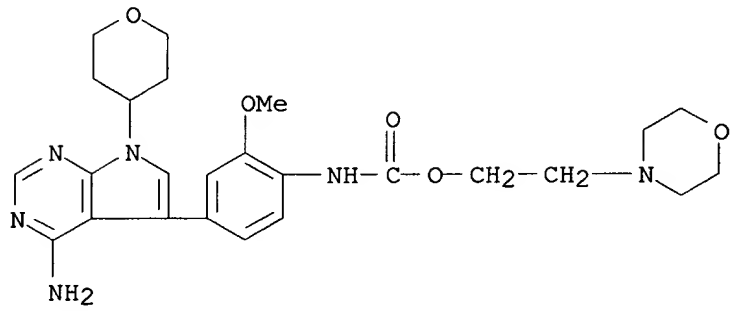
RN 262439-90-7 CAPLUS
CN Carbamic acid, [4-[4-amino-7-(tetrahydro-2H-pyran-4-yl)-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-, 3-pyridinylmethyl ester, monohydrochloride (9CI) (CA INDEX NAME)



● HCl

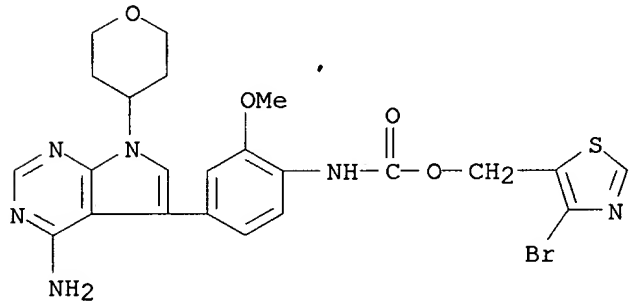
RN 262439-91-8 CAPLUS
CN Carbamic acid, [4-[4-amino-7-(tetrahydro-2H-pyran-4-yl)-7H-pyrrolo[2,3-

d]pyrimidin-5-yl]-2-methoxyphenyl]-, 2-(4-morpholinyl)ethyl ester, monohydrochloride (9CI) (CA INDEX NAME)

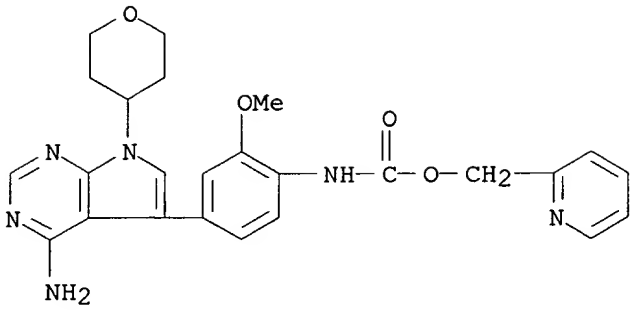


● HCl

RN 262439-92-9 CAPLUS
CN Carbamic acid, [4-[4-amino-7-(tetrahydro-2H-pyran-4-yl)-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-, (4-bromo-5-thiazolyl)methyl ester (9CI) (CA INDEX NAME)

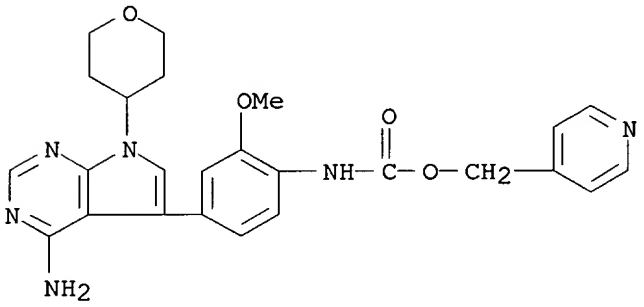


RN 262439-96-3 CAPLUS
CN Carbamic acid, [4-[4-amino-7-(tetrahydro-2H-pyran-4-yl)-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-, 2-pyridinylmethyl ester, monohydrochloride (9CI) (CA INDEX NAME)



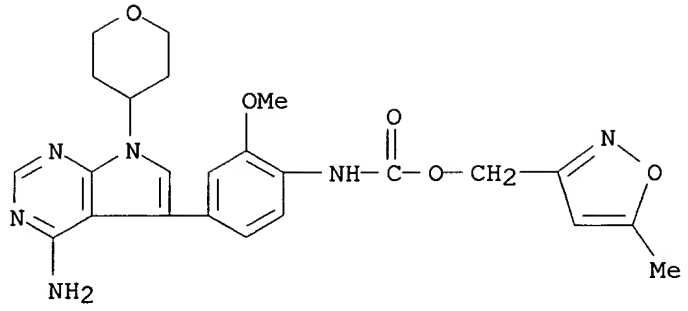
● HCl

RN 262439-97-4 CAPLUS
CN Carbamic acid, [4-[4-amino-7-(tetrahydro-2H-pyran-4-yl)-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-, 4-pyridinylmethyl ester, monohydrochloride (9CI) (CA INDEX NAME)



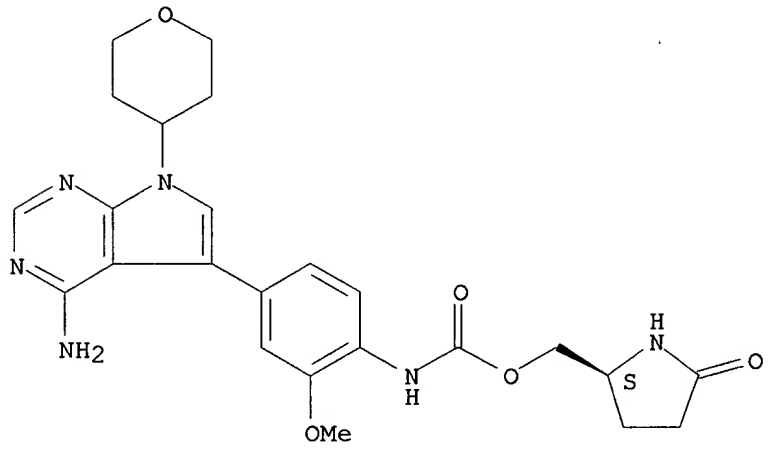
● HCl

RN 262439-98-5 CAPLUS
CN Carbamic acid, [4-[4-amino-7-(tetrahydro-2H-pyran-4-yl)-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-, (5-methyl-3-isoxazolyl)methyl ester (9CI) (CA INDEX NAME)

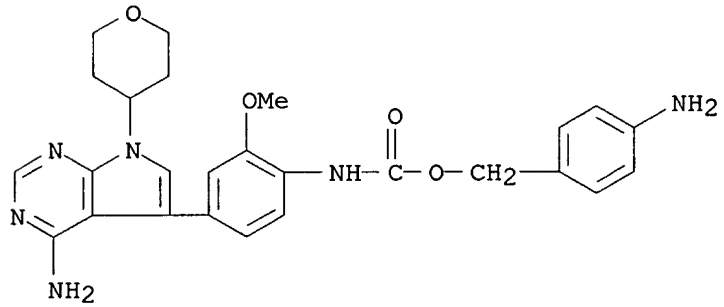


RN 262439-99-6 CAPLUS
CN Carbamic acid, [4-[4-amino-7-(tetrahydro-2H-pyran-4-yl)-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-, [(2S)-5-oxo-2-pyrrolidinyl]methyl ester (9CI) (CA INDEX NAME)

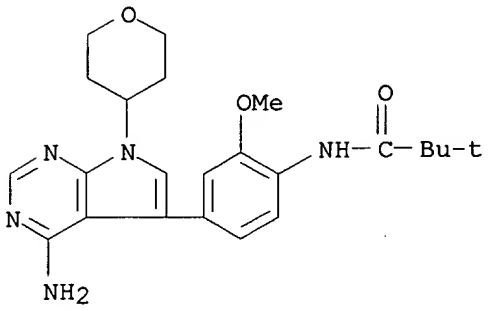
Absolute stereochemistry.



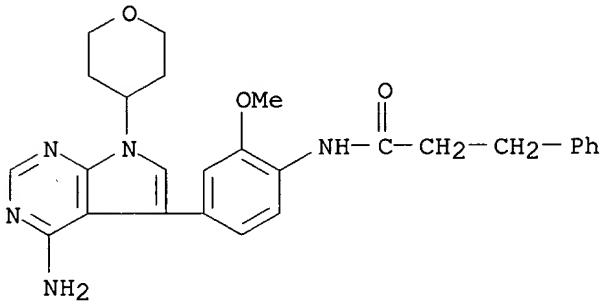
RN 262440-00-6 CAPLUS
CN Carbamic acid, [4-[4-amino-7-(tetrahydro-2H-pyran-4-yl)-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-, (4-aminophenyl)methyl ester (9CI) (CA INDEX NAME)



RN 262440-04-0 CAPLUS
CN Propanamide, N-[4-[4-amino-7-(tetrahydro-2H-pyran-4-yl)-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-2,2-dimethyl- (9CI) (CA INDEX NAME)

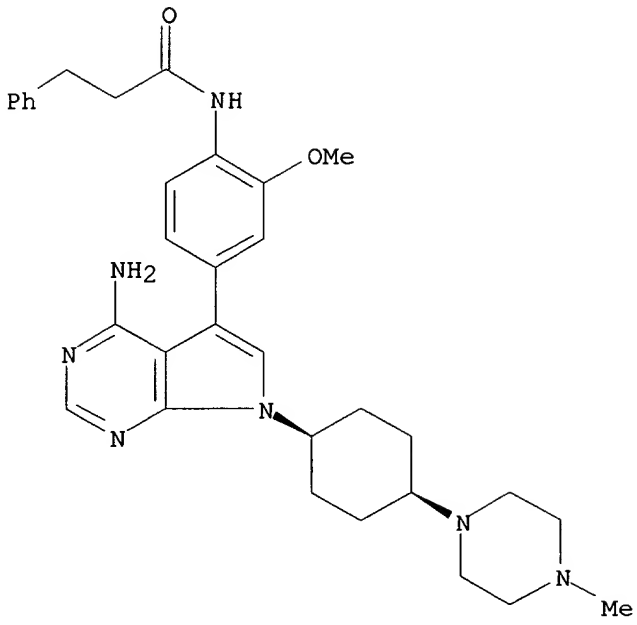


RN 262440-06-2 CAPLUS
CN Benzenepropanamide, N-[4-[4-amino-7-(tetrahydro-2H-pyran-4-yl)-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]- (9CI) (CA INDEX NAME)



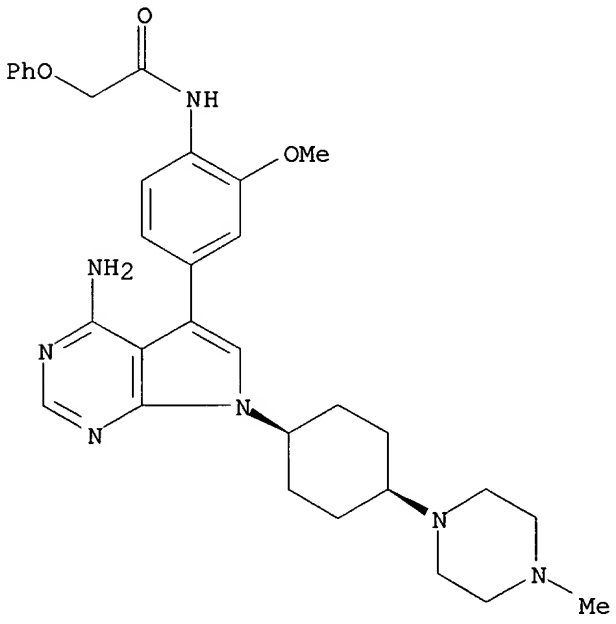
RN 262442-22-8 CAPLUS
CN Benzenepropanamide, N-[4-[4-amino-7-[cis-4-(4-methyl-1-piperazinyl)cyclohexyl]-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]- (9CI) (CA INDEX NAME)

Relative stereochemistry.



RN 262442-32-0 CAPLUS
CN Acetamide, N-[4-[4-amino-7-[cis-4-(4-methyl-1-piperazinyl)cyclohexyl]-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-2-phenoxy- (9CI) (CA INDEX NAME)

Relative stereochemistry.



RN 262442-33-1 CAPLUS
CN Acetamide, N-[4-[4-amino-7-[cis-4-(4-methyl-1-piperazinyl)cyclohexyl]-7H-

pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-2-(4-chlorophenoxy)- (9CI) (CA INDEX NAME)

Relative stereochemistry.

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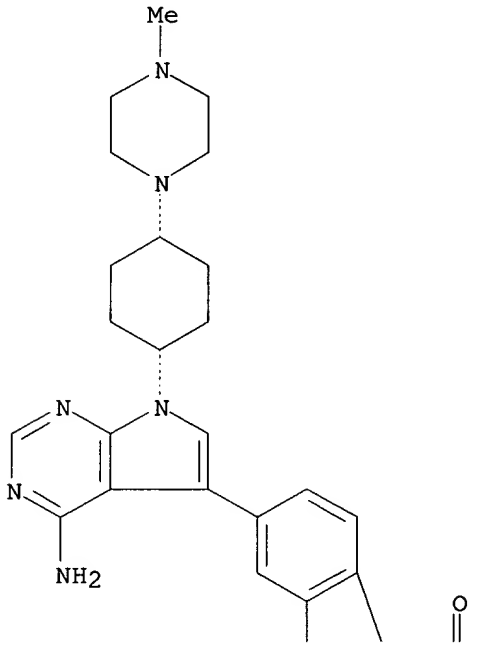

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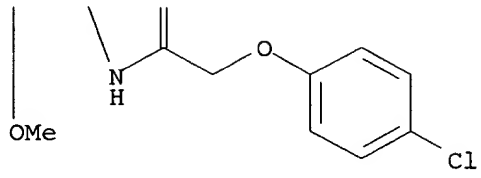
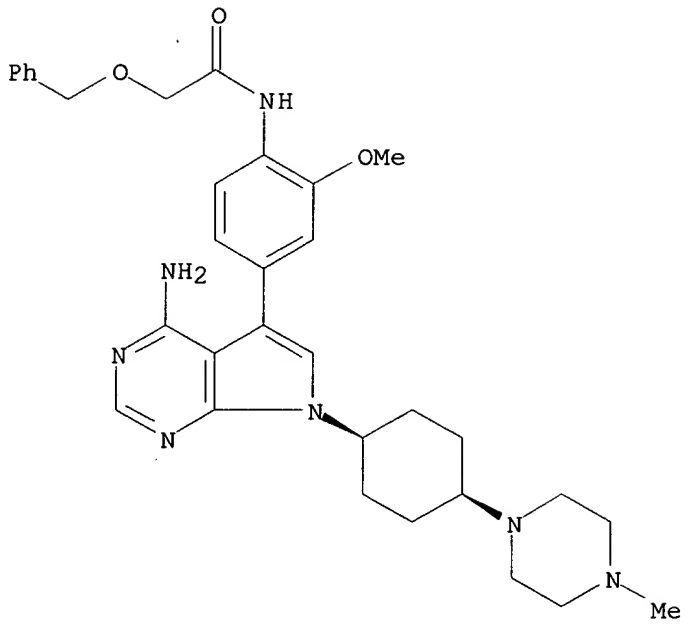


RN 262442-39-7 CAPLUS

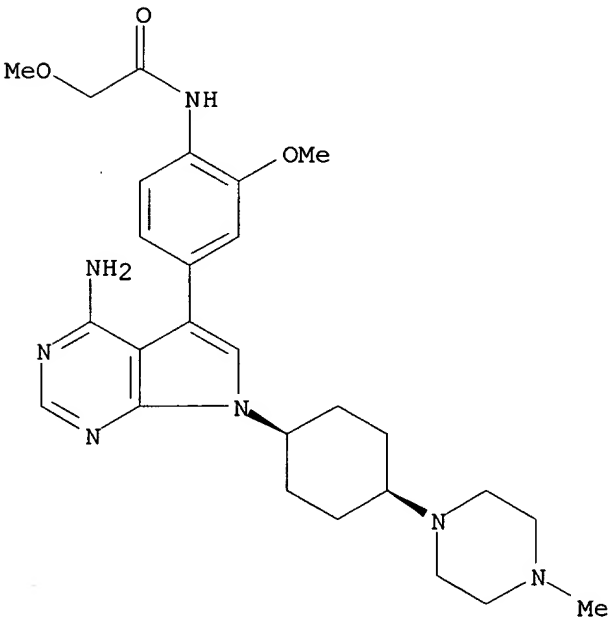
CN Acetamide, N-[4-[4-amino-7-[cis-4-(4-methyl-1-piperazinyl)cyclohexyl]-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-2-(phenylmethoxy)- (9CI) (CA INDEX NAME)

Relative stereochemistry.



RN 262442-47-7 CAPLUS
CN Acetamide, N-[4-[4-amino-7-[cis-4-(4-methyl-1-piperazinyl)cyclohexyl]-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-2-methoxy- (9CI) (CA INDEX NAME)

Relative stereochemistry.



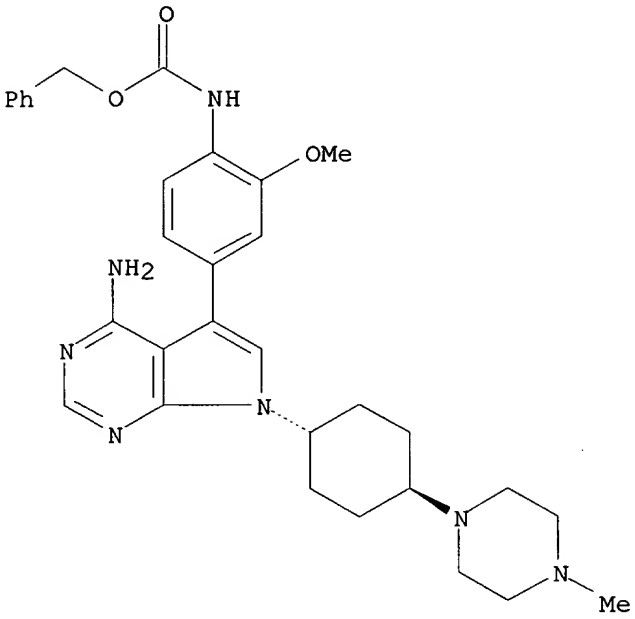
RN 262442-80-8 CAPLUS
CN Carbamic acid, [4-[4-amino-7-[trans-4-(4-methyl-1-piperazinyl)cyclohexyl]-

7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-, phenylmethyl ester, (2Z)-2-butenedioate (1:3) (9CI) (CA INDEX NAME)

CM 1

CRN 262442-79-5
CMF C32 H39 N7 O3

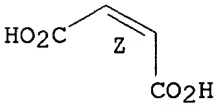
Relative stereochemistry.



CM 2

CRN 110-16-7
CMF C4 H4 O4
CDES 2:Z

Double bond geometry as shown.



RN 262443-69-6 CAPLUS
CN Benzenemethanesulfonamide, N-[4-[4-amino-7-[cis-4-(4-methyl-1-piperazinyl)cyclohexyl]-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-fluorophenyl]-2-nitro-, (2Z)-2-butenedioate (1:3) (9CI) (CA INDEX NAME)

CM 1

CRN 262443-68-5
CMF C30 H35 F N8 O4 S

Relative stereochemistry.

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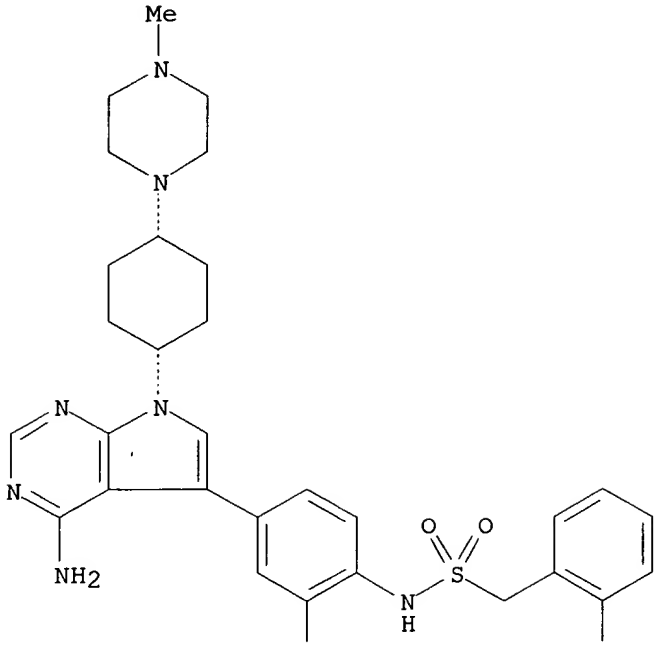

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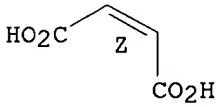
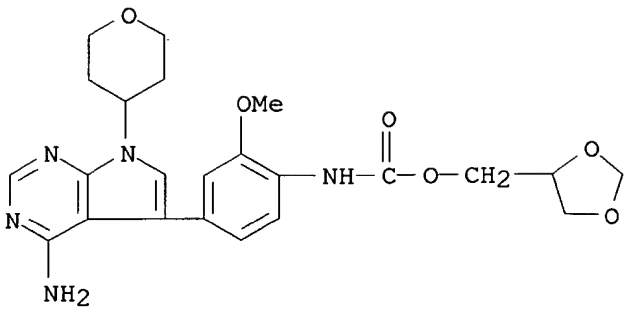
CM 2

CRN 110-16-7
CMF C4 H4 O4
CDES 2:Z

Double bond geometry as shown.



RN 262443-91-4 CAPLUS
CN Carbamic acid, [4-[4-amino-7-(tetrahydro-2H-pyran-4-yl)-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-, 1,3-dioxolan-4-ylmethyl ester (9CI) (CA INDEX NAME)

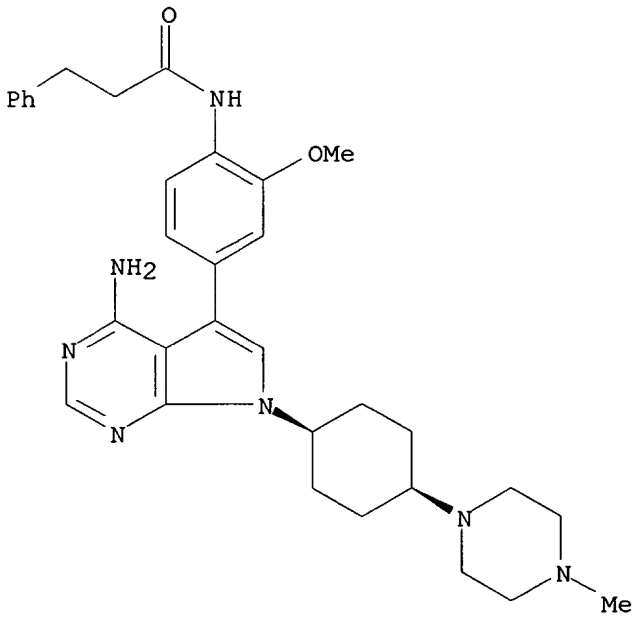


RN 262444-01-9 CAPLUS
CN Benzenepropanamide, N-[4-[4-amino-7-[cis-4-(4-methyl-1-piperazinyl)cyclohexyl]-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-, (2Z)-2-butenedioate (1:3) (9CI) (CA INDEX NAME)

CM 1

CRN 262442-22-8
CMF C33 H41 N7 O2

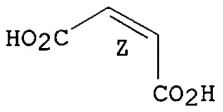
Relative stereochemistry.



CM 2

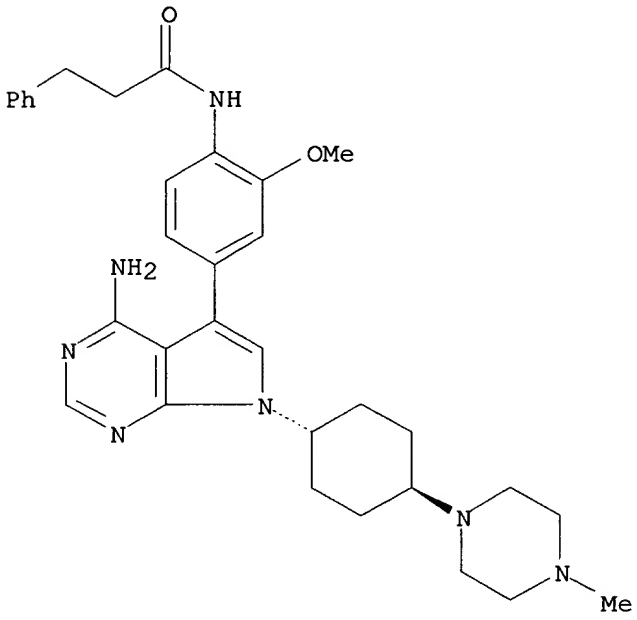
CRN 110-16-7
CMF C4 H4 O4
CDES 2:Z

Double bond geometry as shown.



RN 262444-02-0 CAPLUS
CN Benzenepropanamide, N-[4-[4-amino-7-[trans-4-(4-methyl-1-piperazinyl)cyclohexyl]-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-(9CI) (CA INDEX NAME)

Relative stereochemistry.

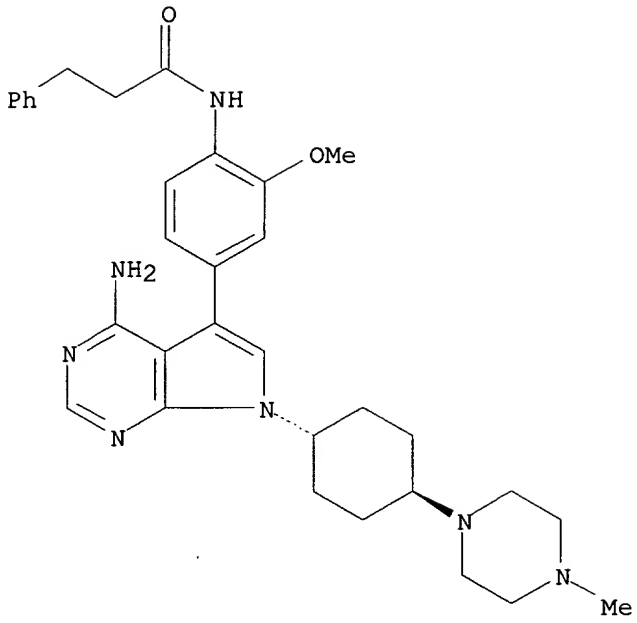


RN 262444-03-1 CAPLUS
CN Benzenepropanamide, N-[4-[4-amino-7-[trans-4-(4-methyl-1-piperazinyl)cyclohexyl]-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-, (2Z)-2-butenedioate (1:3) (9CI) (CA INDEX NAME)

CM 1

CRN 262444-02-0
CMF C33 H41 N7 O2

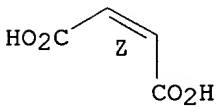
Relative stereochemistry.



CM 2

CRN 110-16-7
CMF C4 H4 O4
CDES 2:Z

Double bond geometry as shown.

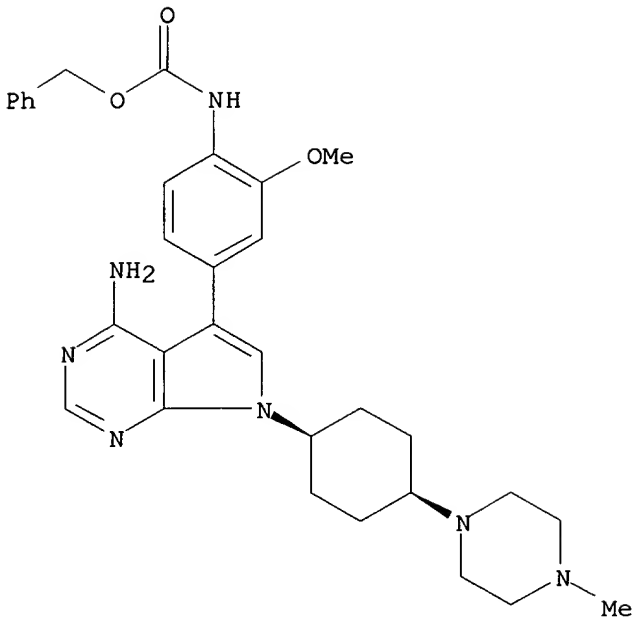


RN 262444-68-8 CAPLUS
CN Carbamic acid, [4-[4-amino-7-[cis-4-(4-methyl-1-piperazinyl)cyclohexyl]-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxyphenyl]-, phenylmethyl ester, (2Z)-2-butenedioate (1:3) (9CI) (CA INDEX NAME)

CM 1

CRN 262444-67-7
CMF C32 H39 N7 O3

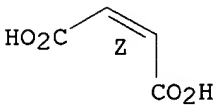
Relative stereochemistry.



CM 2

CRN 110-16-7
CMF C4 H4 O4
CDES 2:Z

Double bond geometry as shown.



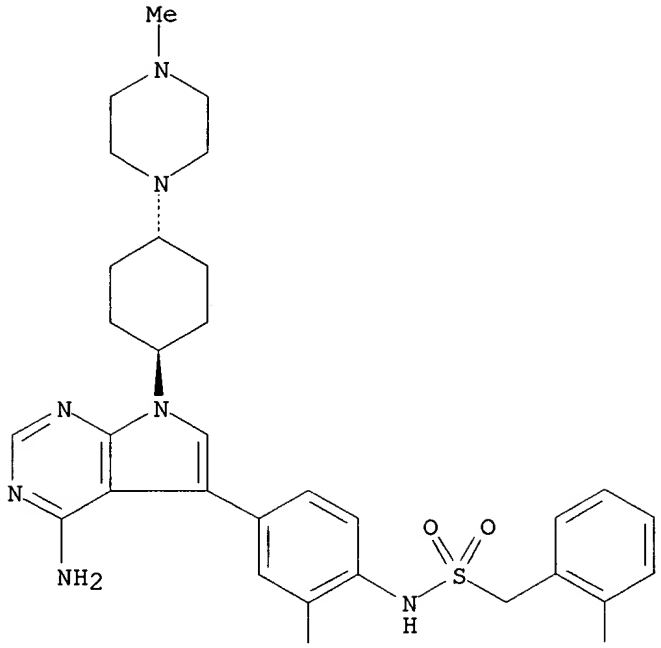
RN 262445-41-0 CAPLUS
CN Benzenemethanesulfonamide, N-[4-[4-amino-7-[trans-4-(4-methyl-1-piperazinyl)cyclohexyl]-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-fluorophenyl]-2-nitro-, (2Z)-2-butenedioate (1:3) (9CI) (CA INDEX NAME)

CM 1

CRN 262445-40-9
CMF C30 H35 F N8 O4 S

Relative stereochemistry.

PAGE 1-A

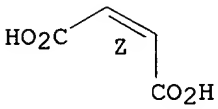


PAGE 2-A

CM 2

CRN 110-16-7
CMF C4 H4 O4
CDES 2:Z

Double bond geometry as shown.



RE.CNT 12 THERE ARE 12 CITED REFERENCES AVAILABLE FOR THIS RECORD
ALL CITATIONS AVAILABLE IN THE RE FORMAT

L4 ANSWER 3 OF 3 CAPLUS COPYRIGHT 2002 ACS
AN 2000:210171 CAPLUS
DN 132:251159
TI Preparation of 4-aminopyrrolopyrimidines as protein kinase inhibitors
IN Calderwood, David; Arnold, Lee D.; Mazdiyasni, Hormoz; Hirst, Gavin; Deng, Bojuan B.
PA BASF Aktiengesellschaft, Germany
SO PCT Int. Appl., 242 pp.
CODEN: PIXXD2
DT Patent
LA English
FAN.CNT 1

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI WO 2000017202	A1	20000330	WO 1999-US21536	19990917
W: AE, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, CA, CH, CN, CR, CU, CZ, DE, DK, DM, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MD, MG, MK, MN, MW, MX, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG, US, UZ, VN, YU, ZA, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM				
RW: GH, GM, KE, LS, MW, SD, SL, SZ, TZ, UG, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG				
AU 9960475	A1	20000410	AU 1999-60475	19990917
EP 1114052	A1	20010711	EP 1999-969414	19990917
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO				
BR 9913888	A	20020108	BR 1999-13888	19990917
NO 2001001357	A	20010514	NO 2001-1357	20010316
PRAI US 1998-100954P	P	19980918		
WO 1999-US21536	W	19990917		

OS MARPAT 132:251159
AB 7H-Pyrrolo[2,3-d]pyrimidin-4-amines (I) [wherein A = (un)substituted 6-membered arom. ring or 5- or 6-membered heteroarom. ring; L = RbN(R)S(O)2, RbN(R)P(O), or RbN(R)P(O)O, where Rb = alkylene group which when taken together with the sulfonamide, phosphinamide or phosphonamide group to which it is bound forms a 5- or 6-membered ring fused to ring A, or L = O, S, N(R), 5-, 6-, or 7-membered (oxa)azaphosphaarom. or (oxa)azaphosphacycloalkyl ring, or a variety of linkers contg. functional groups; R = H, acyl, or (un)substituted aliph., (hetero)arom., or cycloalkyl; R1 = H, 2-Ph-1,3-dioxan-5-yl or (un)substituted (cyclo)alkyl, cycloalkenyl, or phenylalkyl; R2 = H, halo, OH, CN, (un)substituted aliph., cycloalkyl, (hetero)arom., (hetero)aralkyl, amino, or amido; R3 (un)substituted aliph., alkenyl, (hetero)cycloalkyl, or (hetero)arom.; n = 0-6], and physiol. acceptable salts and metabolites thereof, were prepd. For example, II was prepd. in a 6-step sequence involving: (1) amine protection of 4-bromo-2-methoxyaniline with di-tert-Bu dicarbonate, (2) 4-addn. of diboron pinacol ester, (3) 4-substitution with 4-chloro-7-cyclopentyl-5-iodo-7H-pyrrolo[2,3-d]pyrimidine, (4) deprotection of the amine with F3CCO2H, (5) 4-amination of the pyrrolopyrimidine, and (6) addn. of 4-cyanobenzenesulfonyl chloride to the anilino amine. I inhibit serine/threonine and tyrosine kinase activity, affecting immunol., hyperproliferative, and angiogenic processes. All exemplified compds. significantly inhibited either FGFR, PDGFR, KDR, Tie-2, Lck, Fyn, Blk, Lyn, or Src at concns. of .ltoreq. 50 .mu.M, and some significantly inhibited cdc2 at concns. of 50 .ltoreq. .mu.M. Thus, these compds. are useful in the treatment of cancer and hyperproliferative

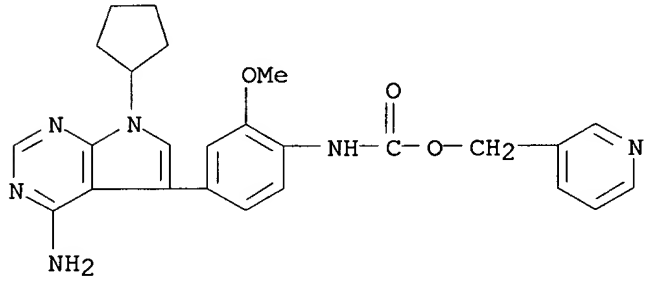
Applicant's PCT

disorders, rheumatoid arthritis, disorders of the immune system, transplant rejections, and inflammatory disorders.

IT **262433-21-6P**
RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); RCT (Reactant); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); RACT (Reactant or reagent); USES (Uses)
(intermediate; prepn. of 7H-pyrrolo[2,3-d]pyrimidin-4-amines as protein kinase inhibitors)

RN 262433-21-6 CAPLUS

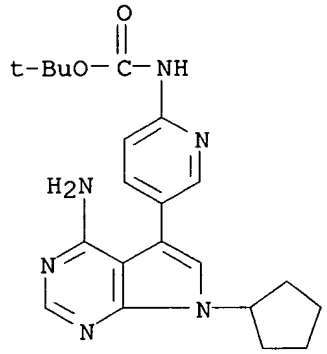
CN Carbamic acid, [4-(4-amino-7-cyclopentyl-7H-pyrrolo[2,3-d]pyrimidin-5-yl)-2-methoxyphenyl]-, 3-pyridinylmethyl ester (9CI) (CA INDEX NAME)



IT **262433-14-7P**
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)
(intermediate; prepn. of 7H-pyrrolo[2,3-d]pyrimidin-4-amines as protein kinase inhibitors)

RN 262433-14-7 CAPLUS

CN Carbamic acid, [5-(4-amino-7-cyclopentyl-7H-pyrrolo[2,3-d]pyrimidin-5-yl)-2-pyridinyl]-, 1,1-dimethylethyl ester (9CI) (CA INDEX NAME)

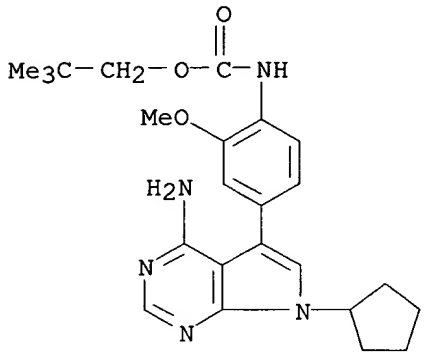


IT **262430-19-3P 262430-20-6P 262430-23-9P**
262430-24-0P 262430-25-1P 262430-26-2P
262430-27-3P 262430-28-4P 262430-29-5P
262430-30-8P 262430-31-9P 262430-32-0P
262430-33-1P 262430-34-2P 262430-35-3P
262432-67-7P 262432-68-8P 262432-69-9P
262433-00-1P
RL: BAC (Biological activity or effector, except adverse); BSU (Biological

study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
(target compd.; prepn. of 7H-pyrrolo[2,3-d]pyrimidin-4-amines as protein kinase inhibitors)

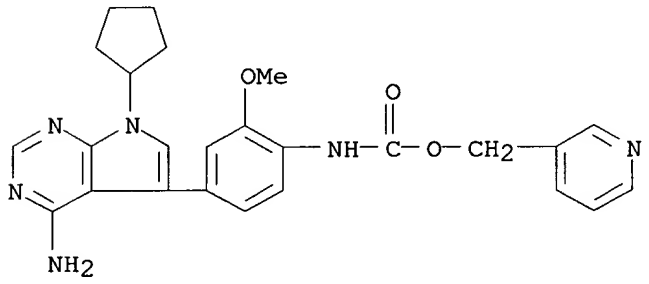
RN 262430-19-3 CAPLUS

CN Carbamic acid, [4-(4-amino-7-cyclopentyl-7H-pyrrolo[2,3-d]pyrimidin-5-yl)-2-methoxyphenyl]-, 2,2-dimethylpropyl ester (9CI) (CA INDEX NAME)



RN 262430-20-6 CAPLUS

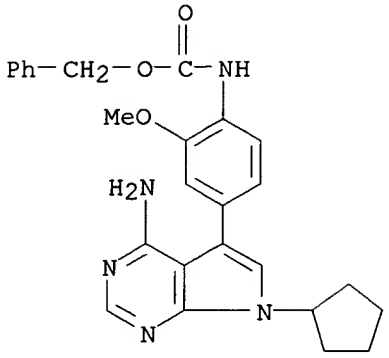
CN Carbamic acid, [4-(4-amino-7-cyclopentyl-7H-pyrrolo[2,3-d]pyrimidin-5-yl)-2-methoxyphenyl]-, 3-pyridinylmethyl ester, hydrochloride (9CI) (CA INDEX NAME)



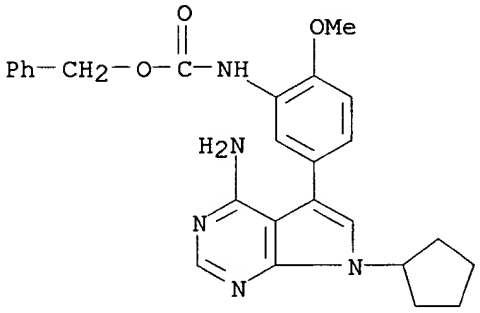
●x HCl

RN 262430-23-9 CAPLUS

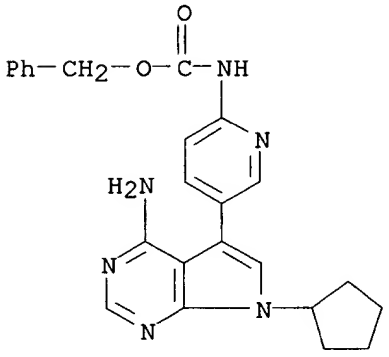
CN Carbamic acid, [4-(4-amino-7-cyclopentyl-7H-pyrrolo[2,3-d]pyrimidin-5-yl)-2-methoxyphenyl]-, phenylmethyl ester (9CI) (CA INDEX NAME)



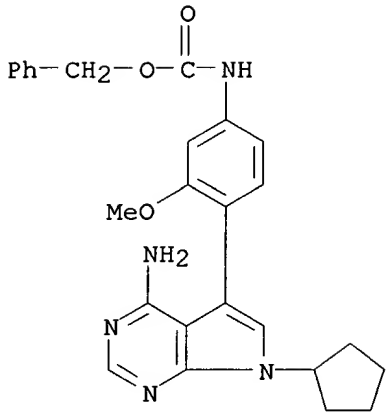
RN 262430-24-0 CAPLUS
CN Carbamic acid, [5-(4-amino-7-cyclopentyl-7H-pyrrolo[2,3-d]pyrimidin-5-yl)-2-methoxyphenyl]-, phenylmethyl ester (9CI) (CA INDEX NAME)



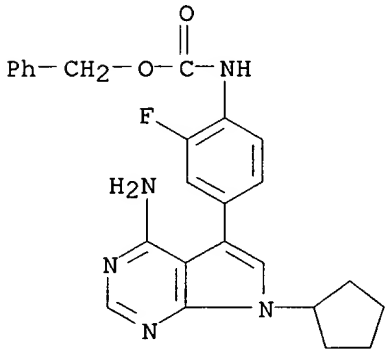
RN 262430-25-1 CAPLUS
CN Carbamic acid, [5-(4-amino-7-cyclopentyl-7H-pyrrolo[2,3-d]pyrimidin-5-yl)-2-pyridinyl]-, phenylmethyl ester (9CI) (CA INDEX NAME)



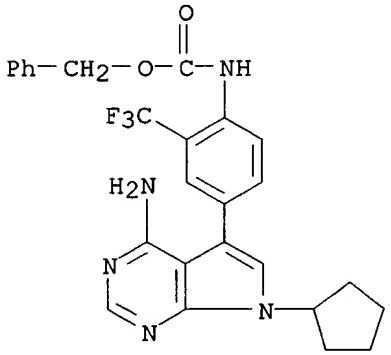
RN 262430-26-2 CAPLUS
CN Carbamic acid, [4-(4-amino-7-cyclopentyl-7H-pyrrolo[2,3-d]pyrimidin-5-yl)-3-methoxyphenyl]-, phenylmethyl ester (9CI) (CA INDEX NAME)



RN 262430-27-3 CAPLUS
CN Carbamic acid, [4-(4-amino-7-cyclopentyl-7H-pyrrolo[2,3-d]pyrimidin-5-yl)-2-fluorophenyl]-, phenylmethyl ester (9CI) (CA INDEX NAME)

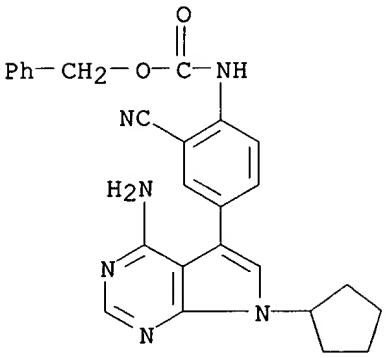


RN 262430-28-4 CAPLUS
CN Carbamic acid, [4-(4-amino-7-cyclopentyl-7H-pyrrolo[2,3-d]pyrimidin-5-yl)-2-(trifluoromethyl)phenyl]-, phenylmethyl ester (9CI) (CA INDEX NAME)

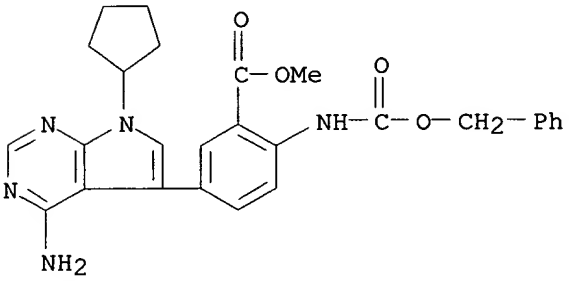


RN 262430-29-5 CAPLUS
CN Carbamic acid, [4-(4-amino-7-cyclopentyl-7H-pyrrolo[2,3-d]pyrimidin-5-yl)-

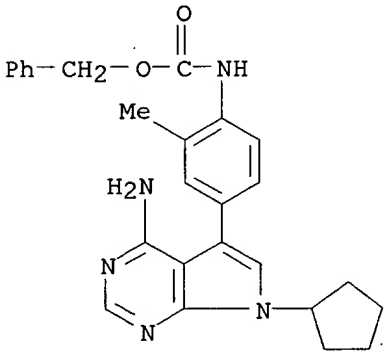
2-cyanophenyl]-, phenylmethyl ester (9CI) (CA INDEX NAME)



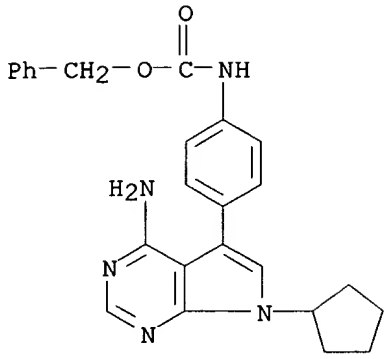
RN 262430-30-8 CAPLUS
CN Benzoic acid, 5-(4-amino-7-cyclopentyl-7H-pyrrolo[2,3-d]pyrimidin-5-yl)-2-[[(phenylmethoxy)carbonyl]amino]-, methyl ester (9CI) (CA INDEX NAME)



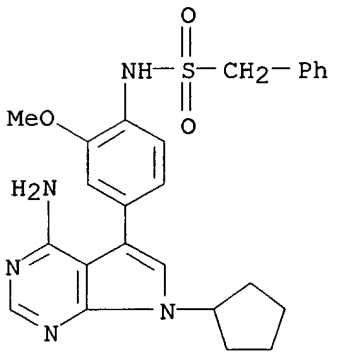
RN 262430-31-9 CAPLUS
CN Carbamic acid, [4-(4-amino-7-cyclopentyl-7H-pyrrolo[2,3-d]pyrimidin-5-yl)-2-methylphenyl]-, phenylmethyl ester (9CI) (CA INDEX NAME)



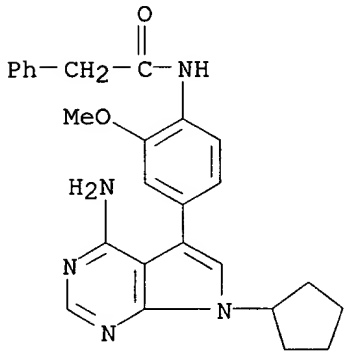
RN 262430-32-0 CAPLUS
CN Carbamic acid, [4-(4-amino-7-cyclopentyl-7H-pyrrolo[2,3-d]pyrimidin-5-yl)phenyl]-, phenylmethyl ester (9CI) (CA INDEX NAME)



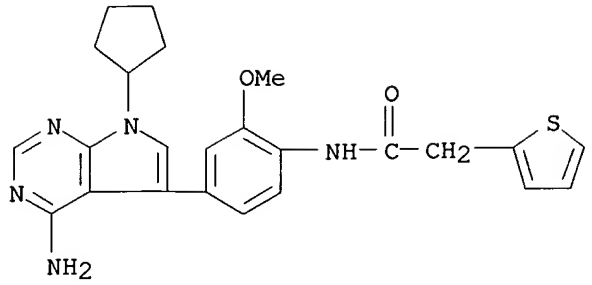
RN 262430-33-1 CAPLUS
CN Benzenemethanesulfonamide, N-[4-(4-amino-7-cyclopentyl-7H-pyrrolo[2,3-d]pyrimidin-5-yl)-2-methoxyphenyl]- (9CI) (CA INDEX NAME)



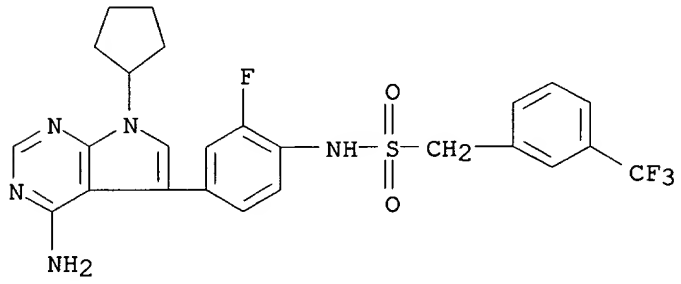
RN 262430-34-2 CAPLUS
CN Benzeneacetamide, N-[4-(4-amino-7-cyclopentyl-7H-pyrrolo[2,3-d]pyrimidin-5-yl)-2-methoxyphenyl]- (9CI) (CA INDEX NAME)



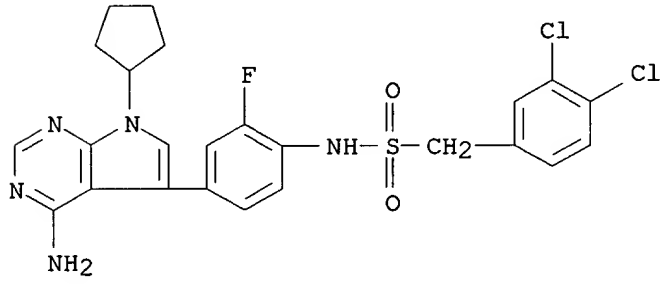
RN 262430-35-3 CAPLUS
CN 2-Thiopheneacetamide, N-[4-(4-amino-7-cyclopentyl-7H-pyrrolo[2,3-d]pyrimidin-5-yl)-2-methoxyphenyl]- (9CI) (CA INDEX NAME)



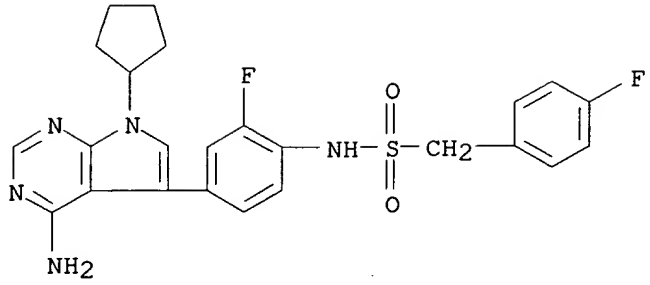
RN 262432-67-7 CAPLUS
CN Benzenemethanesulfonamide, N-[4-(4-amino-7-cyclopentyl-7H-pyrrolo[2,3-d]pyrimidin-5-yl)-2-fluorophenyl]-3-(trifluoromethyl)- (9CI) (CA INDEX NAME)



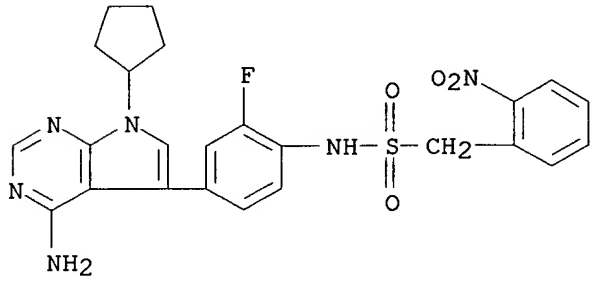
RN 262432-68-8 CAPLUS
CN Benzenemethanesulfonamide, N-[4-(4-amino-7-cyclopentyl-7H-pyrrolo[2,3-d]pyrimidin-5-yl)-2-fluorophenyl]-3,4-dichloro- (9CI) (CA INDEX NAME)



RN 262432-69-9 CAPLUS
CN Benzenemethanesulfonamide, N-[4-(4-amino-7-cyclopentyl-7H-pyrrolo[2,3-d]pyrimidin-5-yl)-2-fluorophenyl]-4-fluoro- (9CI) (CA INDEX NAME)



RN 262433-00-1 CAPLUS
CN Benzenemethanesulfonamide, N-[4-(4-amino-7-cyclopentyl-7H-pyrrolo[2,3-d]pyrimidin-5-yl)-2-fluorophenyl]-2-nitro- (9CI) (CA INDEX NAME)



RE.CNT 6 THERE ARE 6 CITED REFERENCES AVAILABLE FOR THIS RECORD
ALL CITATIONS AVAILABLE IN THE RE FORMAT

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     FILE 'REGISTRY' ENTERED AT 11:32:24 ON 17 MAY 2002
L1              STRUCTURE UPLOADED
L2            4 S L1
L3           61 S L1 SSS FUL

     FILE 'CAPLUS' ENTERED AT 11:34:00 ON 17 MAY 2002
L4            3 S L3

     FILE 'CAOLD' ENTERED AT 11:34:42 ON 17 MAY 2002

=> s l3
L5            0 L3

=> log y
COST IN U.S. DOLLARS                                SINCE FILE      TOTAL
                                                       ENTRY    SESSION
FULL ESTIMATED COST                                     0.38     155.19

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)          SINCE FILE      TOTAL
                                                       ENTRY    SESSION
CA SUBSCRIBER PRICE                                     0.00      -1.86

STN INTERNATIONAL LOGOFF AT 11:34:56 ON 17 MAY 2002
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L1 HAS NO ANSWERS
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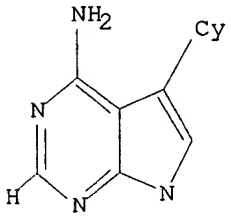


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Structure attributes must be viewed using STN Express query preparation.


=> s l1 sss sam

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100.0% PROCESSED     319 ITERATIONS                          21 ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS:  ONLINE  **COMPLETE**
                        BATCH   **COMPLETE**
PROJECTED ITERATIONS:         5309 TO     7451
PROJECTED ANSWERS:             146 TO      694

L2           21 SEA SSS SAM L1

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Uploading 09399083.str

L3      STRUCTURE UPLOADED

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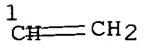

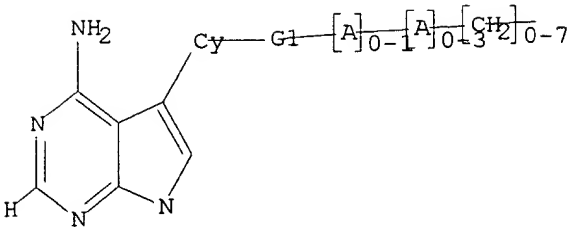

G1 O,S,N,CH2,CH,[@1-@2]

Structure attributes must be viewed using STN Express query preparation.

=> s l3 sss sam

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100.0% PROCESSED 318 ITERATIONS 7 ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 5291 TO 7429
PROJECTED ANSWERS: 7 TO 298

L4 7 SEA SSS SAM L3

=> d his

(FILE 'HOME' ENTERED AT 16:50:35 ON 06 MAR 2000)

FILE 'REGISTRY' ENTERED AT 16:50:52 ON 06 MAR 2000
L1 STRUCTURE UPLOADED
L2 21 S L1 SSS SAM

FILE 'STNGUIDE' ENTERED AT 16:52:10 ON 06 MAR 2000

FILE 'REGISTRY' ENTERED AT 17:00:44 ON 06 MAR 2000
L3 STRUCTURE UPLOADED
L4 7 S L3 SSS SAM

=> s l3 sss ful

FULL SEARCH INITIATED 17:03:39 FILE 'REGISTRY'

FULL SCREEN SEARCH COMPLETED - 6282 TO ITERATE
100.0% PROCESSED 6282 ITERATIONS 142 ANSWERS
SEARCH TIME: 00.00.04

L5 142 SEA SSS FUL L3

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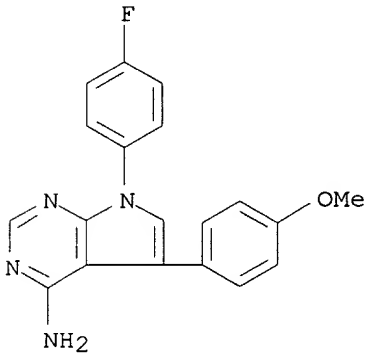
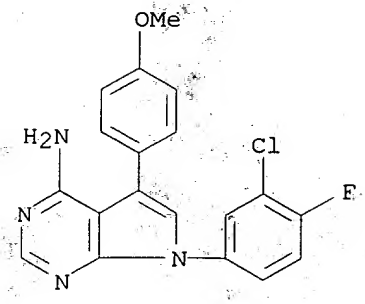
L6 10 L5

=> d l6 1-10 bib,ab,hitstr

L6 ANSWER 1 OF 10 CAPLUS COPYRIGHT 2000 ACS
AN 1999:457818 CAPLUS
DN 131:214251
TI Synthesis and reactions of fluoroaryl substituted 2-amino-3-cyanopyrroles and pyrrolo[2,3-d]pyrimidines
AU Dave, Chaitanya G.; Desai, Nirmal D.
CS Organic Syntheses Laboratory, M. G. Science Institute, Ahmedabad, 380 009, India
SO J. Heterocycl. Chem. (1999), 36(3), 729-733
CODEN: JHTCAD; ISSN: 0022-152X
PB HeteroCorporation
DT Journal
LA English
AB Some fluoroaryl substituted 2-amino-3-cyanopyrroles were synthesized from the reaction between (2-bromo-1-arylalkylidene)propanedinitriles and fluoroaryl substituted arom. amines under Gewald reaction condition, which on reaction with formamide and formic acid gave 4-aminopyrrolo[2,3-d]pyrimidines and pyrrolo[2,3-d]-pyrimidin-4(3H)-ones (4), resp. 4-Chloropyrrolo[2,3-d]pyrimidines were prepd. by chlorination of 4 with P oxychloride, which on hydrazinolysis provided 4-hydrazinopyrrolo[2,3-d]pyrimidines.
IT **243665-89-6P 243665-90-9P**
RL: SPN (Synthetic preparation); PREP (Preparation)
(synthesis and reactions of fluoroaryl substituted aminocyanopyrroles and pyrrolo[2,3-d]pyrimidines)
RN 243665-89-6 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 7-(4-fluorophenyl)-5-(4-methoxyphenyl)- (9CI) (CA INDEX NAME)

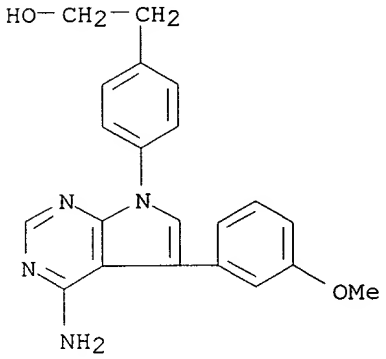


RN 243665-90-9 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 7-(3-chloro-4-fluorophenyl)-5-(4-methoxyphenyl)- (9CI) (CA INDEX NAME)

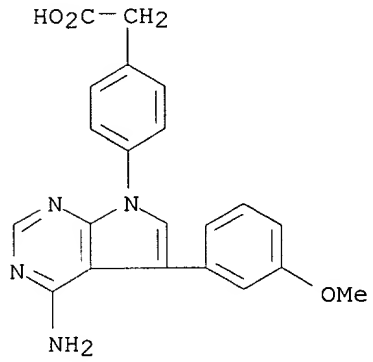

OMe
H2N
Cl
F
N
N
N

L6 ANSWER 2 OF 10 CAPLUS COPYRIGHT 2000 ACS
AN 1999:309002 CAPLUS
DN 131:124921
TI A novel inhibitor of the tyrosine kinase Src suppresses phosphorylation of its major cellular substrates and reduces bone resorption in vitro and in rodent models in vivo
AU Missbach, M.; Jeschke, M.; Feyen, J.; Muller, K.; Glatt, M.; Green, J.; Susa, M.
CS Novartis Pharma AG, Research Bone Metabolism, Basel, CH-4002, Switz.
SO Bone (N. Y.) (1999), 24(5), 437-449
CODEN: BONEDL; ISSN: 8756-3282
PB Elsevier Science Inc.
DT Journal
LA English
AB The tyrosine kinase Src has been implicated in the process of osteoclast-mediated bone resorption. Here, we describe a novel class of Src inhibitors, substituted 5,7-diphenyl-pyrrolo[2,3-d]pyrimidines, and characterize one of them, CGP77675, in vitro and in models of bone resorption in vivo. In vitro, CGP77675 inhibited phosphorylation of peptide substrates and autophosphorylation of purified Src (concn. producing half-maximal inhibition [IC50] values 5-20 and 40 nmol/L, resp.). The compd. was selective toward other protein kinases: the Src IC50 value was lower than those for Cdc2 (>500-fold), epidermal growth factor (EGF) receptor (7.5-fold), and vascular endothelial growth factor receptor (>50-fold), and for v-Abl (15-fold) and focal adhesion kinase (Fak) (>25-fold). The Src kinase family members Lck and Yes were inhibited with IC50 values 20-fold higher than or equal to Src. To measure the inhibition of cellular Src activity, we identified the major tyrosine-phosphorylated proteins in an Src-overexpressing cell line IC8.1 as Src, Fak, and paxillin. CGP77675 potently inhibited tyrosine phosphorylation of the Src substrates Fak and paxillin, but had much less effect on Src (IC50 values 0.3, 0.5, and 5.7 .mu.mol/L). The phosphorylation of Src in IC8.1 cells reflected phosphorylation of the neg. regulatory tyrosine 527 (Y527); thus, the inhibitor was selective against the Y527 C-terminal Src kinase Csk. In osteoblastic MC3T3-E1 cells, CGP77675 inhibited signaling induced by PDGF at the receptor level, but not signaling by EGF, basic fibroblast growth factor, insulin-like growth factor-1, and phorbol 12-myristate 13-acetate. The effect of CGP77675 on bone resorption was evaluated in vitro and in vivo. The parathyroid hormone-induced bone resorption in rat fetal long bone cultures was inhibited with an IC50 of 0.8 .mu.mol/L. CGP77675 dose-dependently reduced the hypercalcemia induced in mice by interleukin-1.beta. and partly prevented bone loss and microarchitectural changes in young ovariectomized rats, showing that the protective effect on bone was exerted via the inhibition of bone resorption. Thus, specific Src family kinase inhibitors may be useful for the treatment of diseases assocd. with elevated bone loss.
IT **234772-58-8 234772-59-9 234772-60-2 234772-61-3 234772-62-4 234772-63-5**
RL: BAC (Biological activity or effector, except adverse); PRP (Properties); BIOL (Biological study)
(inhibitor of tyrosine kinase Src suppresses phosphorylation of its major cellular substrates and reduces bone resorption in vitro and in

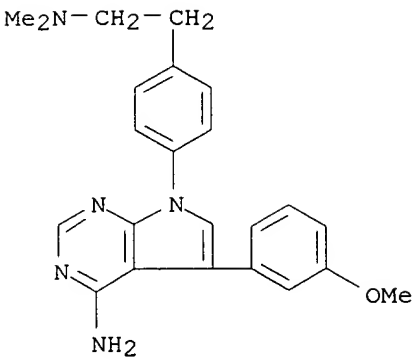
rodent models in vivo)
RN 234772-58-8 CAPLUS
CN Benzeneethanol,
4-[4-amino-5-(3-methoxyphenyl)-7H-pyrrolo[2,3-d]pyrimidin-
7-yl]- (9CI) (CA INDEX NAME)



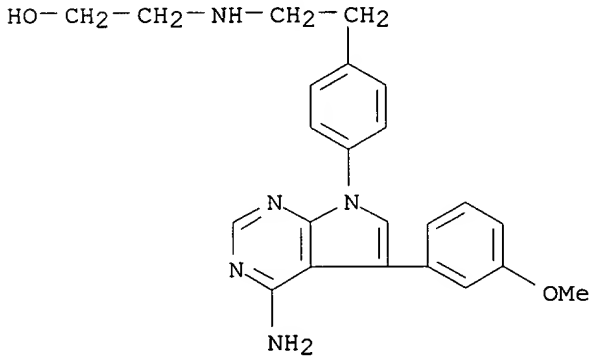
RN 234772-59-9 CAPLUS
CN Benzeneacetic acid, 4-[4-amino-5-(3-methoxyphenyl)-7H-pyrrolo[2,3-
d]pyrimidin-7-yl]- (9CI) (CA INDEX NAME)



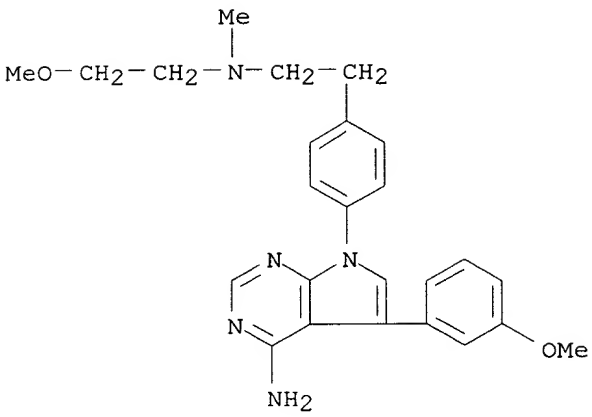
RN 234772-60-2 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine,
7-[4-[2-(dimethylamino)ethyl]phenyl]-5-
(3-methoxyphenyl)- (9CI) (CA INDEX NAME)



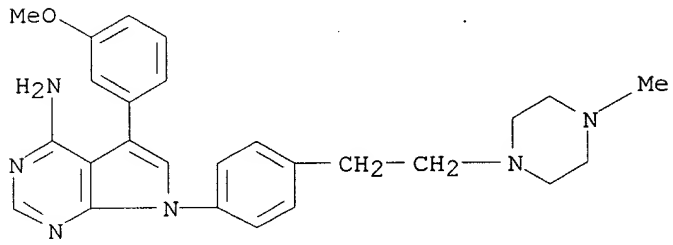
RN 234772-61-3 CAPLUS
CN Ethanol,
2-[[2-[4-[4-amino-5-(3-methoxyphenyl)-7H-pyrrolo[2,3-d]pyrimidin-7-yl]phenyl]ethyl]amino]- (9CI) (CA INDEX NAME)



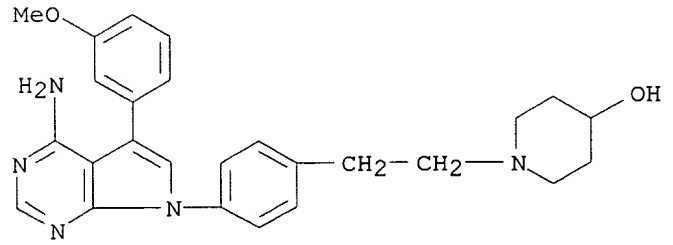
RN 234772-62-4 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 7-[4-[2-[(2-methoxyethyl)methylamino]ethyl]phenyl]-5-(3-methoxyphenyl)- (9CI) (CA INDEX NAME)



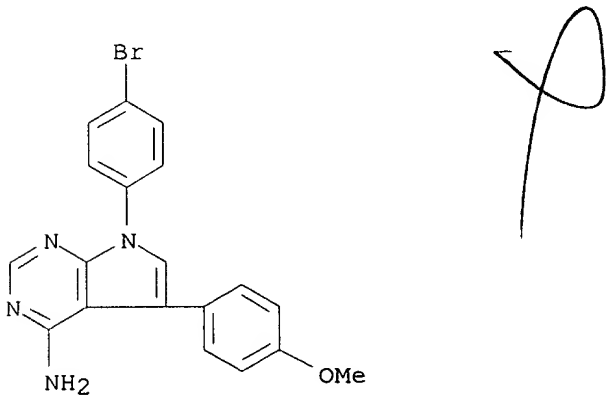
RN 234772-63-5 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 5-(3-methoxyphenyl)-7-[4-[2-(4-methyl-1-piperazinyl)ethyl]phenyl]- (9CI) (CA INDEX NAME)



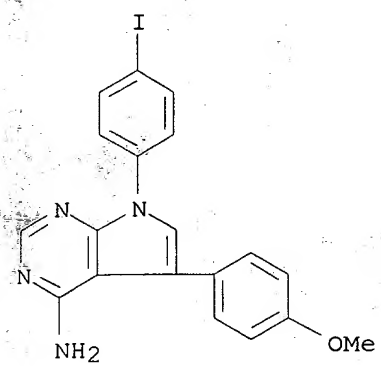
IT **234772-64-6**
RL: BAC (Biological activity or effector, except adverse); PRP (Properties); THU (Therapeutic use); BIOL (Biological study); USES (Uses)
(inhibitor of tyrosine kinase Src suppresses phosphorylation of its major cellular substrates and reduces bone resorption in vitro and in rodent models in vivo)
RN 234772-64-6 CAPLUS
CN 4-Piperidinol, 1-[2-[4-[4-amino-5-(3-methoxyphenyl)-7H-pyrrolo[2,3-d]pyrimidin-7-yl]phenyl]ethyl]- (9CI) (CA INDEX NAME)



L6 ANSWER 3 OF 10 CAPLUS COPYRIGHT 2000 ACS
AN 1999:51036 CAPLUS
DN 130:196617
TI Synthesis of 7H-Tetrazolo[1,5-c]pyrrolo[3,2-e]pyrimidines and their reductive ring cleavage to 4-aminopyrrolo[2,3-d]pyrimidines
AU Dave, Chaitanya G.; Shah, Rina D.
CS Organic Syntheses Laboratory, M. G. Science Institute, Ahmedabad, 300 009, India
SO J. Heterocycl. Chem. (1998), 35(6), 1295-1300
CODEN: JHTCAD; ISSN: 0022-152X
PB HeteroCorporation
DT Journal
LA English
AB Some new 7,9-disubstituted 7H-tetrazolo[1,5-c]pyrrolo[3,2-e]pyrimidines (I; R = H, MeO, Cl; R1 = MeO, Br, I, Me) have been synthesized either by diazotization of 4-hydrazino-7H-pyrrolo[2,3-d]pyrimidines (II; same R, R1; R2 = NHNH2), obtained by hydrazinolysis of II (R2 = Cl) or via a substitution reaction between II (R2 = Cl) and sodium azide. 5,7-Disubstituted 7H-pyrrolo[2,3-d]pyrimidin-4(3H)-ones were obtained by cyclocondensation of 1,4-disubstituted 2-amino-3-cyanopyrroles with formic acid; subsequent chlorination using phosphorus oxychloride afforded II (R2 = Cl). A novel route to II (R2 = NH2) via reductive ring cleavage of I has been reported.
IT **121405-32-1P 121405-33-2P**
RL: SPN (Synthetic preparation); PREP (Preparation)
(prepn. of 7H-tetrazolo[1,5-c]pyrrolo[3,2-e]pyrimidines and their reductive ring cleavage to 4-aminopyrrolo[2,3-d]pyrimidines)
RN 121405-32-1 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 7-(4-bromophenyl)-5-(4-methoxyphenyl)- (9CI) (CA INDEX NAME)



RN 121405-33-2 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 7-(4-iodophenyl)-5-(4-methoxyphenyl)- (9CI) (CA INDEX NAME)



9

L6 ANSWER 4 OF 10 CAPLUS COPYRIGHT 2000 ACS
AN 1998:640260 CAPLUS
DN 129:275922
TI Preparation of pyrrolo[2,3-d]pyrimidines as tyrosine kinase inhibitors
IN Calderwood, David John; Johnston, David Norman; Rafferty, Paul; Twigger, Helen Louise; Munschauer, Rainer; Arnold, Lee
PA Knoll A.-G., Germany
SO PCT Int. Appl., 72 pp.
CODEN: PIXXD2
DT Patent
LA English

Applicant's PCT

FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	WO 9841525	A1	19980924	WO 1998-EP1357	19980309
	W: AL, AU, BG, BR, BY, CA, CN, CZ, GE, HU, ID, IL, JP, KR, KZ, LT, LV, MX, NO, NZ, PL, RO, RU, SG, SI, SK, TR, UA, US, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM				
	RW: AT, BE, CH, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE				
	AU 9868293	A1	19981012	AU 1998-68293	19980309
	EP 970084	A1	20000112	EP 1998-913690	19980309
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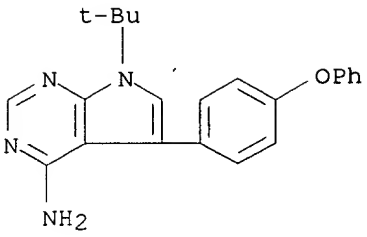
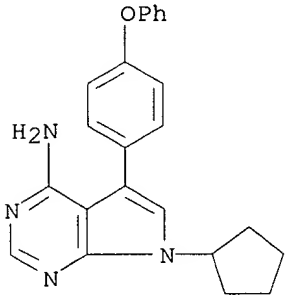
	NO 9904509	A	19990917	NO 1999-4509	19990917
PRAI	US 1997-40836	19970319			
	WO 1998-EP1357	19980309			

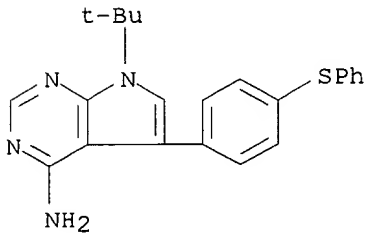
OS MARPAT 129:275922
AB The title compds. [I; R1 = H, 2-phenyl-1,3-dioxan-5-yl, C1-6 alkyl, etc.; R2 = H, C1-6 alkyl, C3-8 cycloalkyl, etc.; R3 = (un)substituted II (A = NH, O, NHSO2, etc.; R5 = (un)substituted Ph, and, addnl., when A is absent
R5 = phthalimido optionally substituted by halo, (un)substituted pyrazolylamino)], useful in treating proliferative diseases and disorders of the immune system in mammals (no data), were prepd. Thus, heating a mixt. of 2-amino-3-cyano-4-(4-phenoxyphenyl)-1-tert-butylpyrrole (prepn. described), formamide and 4-dimethylaminopyridine at 180.degree. for 6 h afforded I [R1 = tBu; R2 = H; R3 = 4-(PhO)C6H4].
IT **213743-29-4P 213743-31-8P 213743-36-3P**
213743-38-5P 213743-44-3P 213743-46-5P
213743-52-3P 213743-74-9P 213743-84-1P
213743-90-9P 213743-94-3P 213744-00-4P
213744-08-2P 213744-10-6P 213744-26-4P
RL: BAC (Biological activity or effector, except adverse); RCT (Reactant);
SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
(prepn. of pyrrolo[2,3-d]pyrimidines as tyrosine kinase inhibitors)
RN 213743-29-4 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 7-(1,1-dimethylethyl)-5-(4-phenoxyphenyl)- (9CI) (CA INDEX NAME)



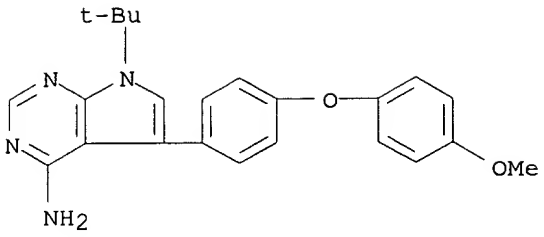
RN 213743-31-8 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 7-cyclopentyl-5-(4-phenoxyphenyl)- (9CI) (CA INDEX NAME)



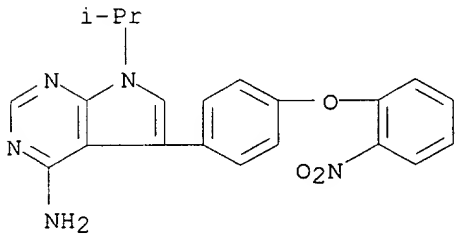
RN 213743-36-3 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 7-(1,1-dimethylethyl)-5-[4-(phenylthio)phenyl]- (9CI) (CA INDEX NAME)



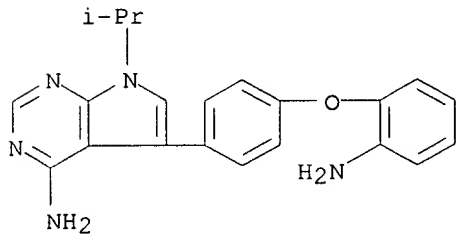
RN 213743-38-5 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 7-(1,1-dimethylethyl)-5-[4-(4-methoxyphenoxy)phenyl]- (9CI) (CA INDEX NAME)



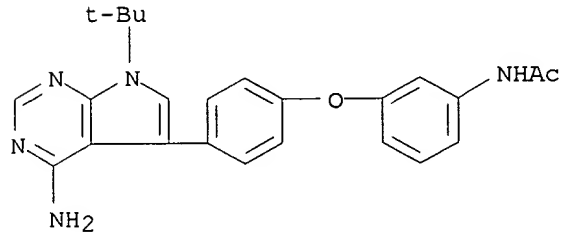
RN 213743-44-3 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 7-(1-methylethyl)-5-[4-(2-nitrophenoxy)phenyl]- (9CI) (CA INDEX NAME)



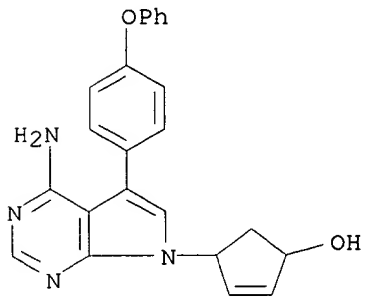
RN 213743-46-5 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 5-[4-(2-aminophenoxy)phenyl]-7-(1-methylethyl)- (9CI) (CA INDEX NAME)



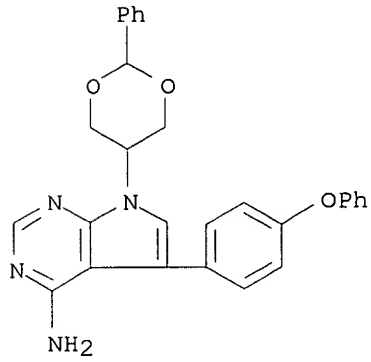
RN 213743-52-3 CAPLUS
CN Acetamide, N-[3-[4-[4-amino-7-(1,1-dimethylethyl)-7H-pyrrolo[2,3-d]pyrimidin-5-yl]phenoxy]phenyl]- (9CI) (CA INDEX NAME)



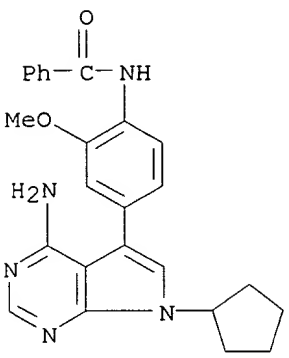
RN 213743-74-9 CAPLUS
CN 2-Cyclopenten-1-ol, 4-[4-amino-5-(4-phenoxyphenyl)-7H-pyrrolo[2,3-d]pyrimidin-7-yl]- (9CI) (CA INDEX NAME)



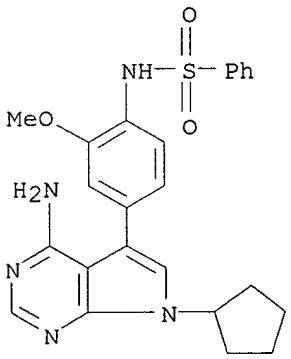
RN 213743-84-1 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 5-(4-phenoxyphenyl)-7-(2-phenyl-1,3-dioxan-5-yl)- (9CI) (CA INDEX NAME)



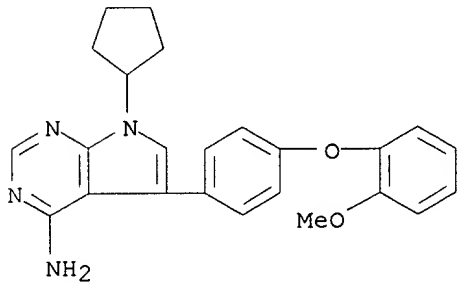
RN 213743-90-9 CAPLUS
CN Benzamide,
N-[4-(4-amino-7-cyclopentyl-7H-pyrrolo[2,3-d]pyrimidin-5-yl)-2-methoxyphenyl]- (9CI) (CA INDEX NAME)



RN 213743-94-3 CAPLUS
CN Benzenesulfonamide,
N-[4-(4-amino-7-cyclopentyl-7H-pyrrolo[2,3-d]pyrimidin-
5-yl)-2-methoxyphenyl]- (9CI) (CA INDEX NAME)

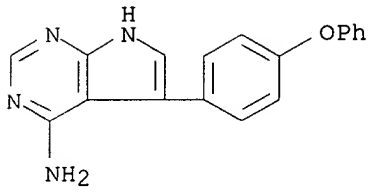


RN 213744-00-4 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 7-cyclopentyl-5-[4-(2-
methoxyphenoxy)phenyl]- (9CI) (CA INDEX NAME)



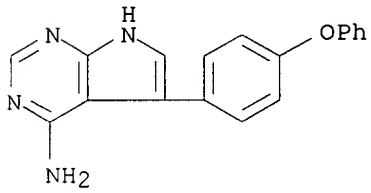
RN 213744-08-2 CAPLUS
CN 1H-Pyrrolo[2,3-d]pyrimidin-4-amine, 5-(4-phenoxyphenyl)-,
monohydrobromide

(9CI) (CA INDEX NAME)

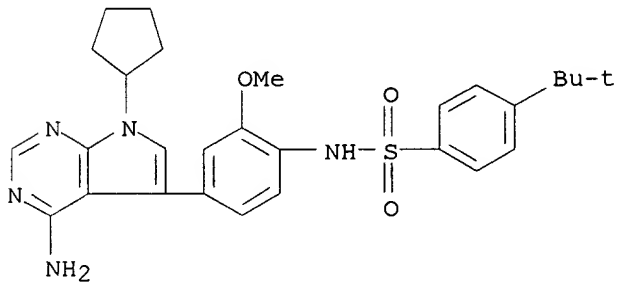


● HBr

RN 213744-10-6 CAPLUS
CN 1H-Pyrrolo[2,3-d]pyrimidin-4-amine, 5-(4-phenoxyphenyl)- (9CI) (CA INDEX NAME)



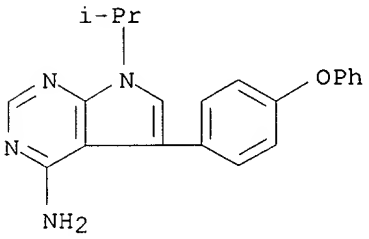
RN 213744-26-4 CAPLUS
CN Benzenesulfonamide,
N-[4-(4-amino-7-cyclopentyl-7H-pyrrolo[2,3-d]pyrimidin-5-yl)-2-methoxyphenyl]-4-(1,1-dimethylethyl)- (9CI) (CA INDEX NAME)



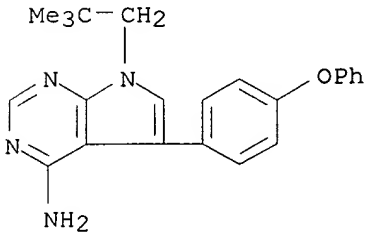
IT **213743-30-7P 213743-34-1P 213743-40-9P**
213743-42-1P 213743-48-7P 213743-50-1P
213743-54-5P 213743-56-7P 213743-58-9P
213743-60-3P 213743-62-5P 213743-64-7P
213743-66-9P 213743-68-1P 213743-70-5P
213743-72-7P 213743-76-1P 213743-78-3P
213743-80-7P 213743-82-9P 213743-86-3P
213743-88-5P 213743-92-1P 213743-96-5P

213743-98-7P 213744-02-6P 213744-04-8P
213744-06-0P 213744-12-8P 213744-14-0P
213744-16-2P 213744-18-4P 213744-20-8P
213744-22-0P 213744-24-2P 213744-29-7P
RL: BAC (Biological activity or effector, except adverse); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
(prepn. of pyrrolo[2,3-d]pyrimidines as tyrosine kinase inhibitors)

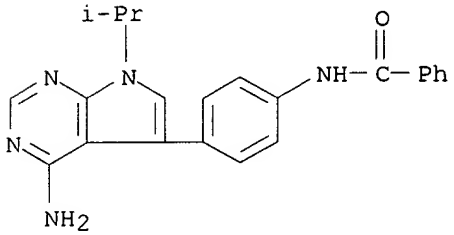
RN 213743-30-7 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 7-(1-methylethyl)-5-(4-phenoxyphenyl)- (9CI) (CA INDEX NAME)



RN 213743-34-1 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 7-(2,2-dimethylpropyl)-5-(4-phenoxyphenyl)- (9CI) (CA INDEX NAME)

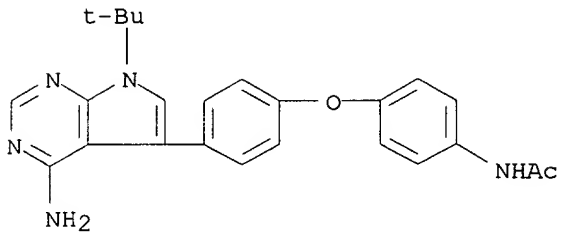


RN 213743-40-9 CAPLUS
CN Benzamide, N-[4-[4-amino-7-(1-methylethyl)-7H-pyrrolo[2,3-d]pyrimidin-5-yl]phenyl]- (9CI) (CA INDEX NAME)

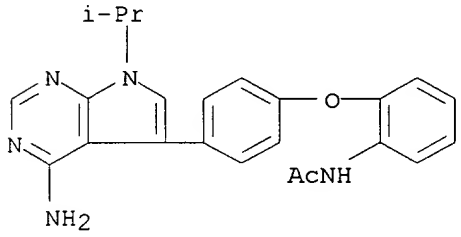


RN 213743-42-1 CAPLUS
CN Acetamide, N-[4-[4-[4-amino-7-(1,1-dimethylethyl)-7H-pyrrolo[2,3-

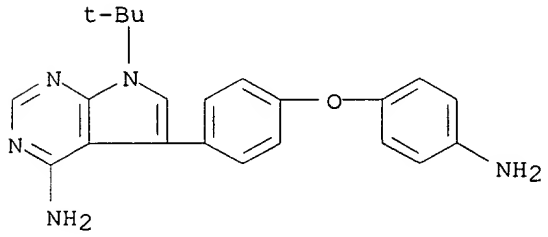
d]pyrimidin-5-yl]phenoxy]phenyl]- (9CI) (CA INDEX NAME)



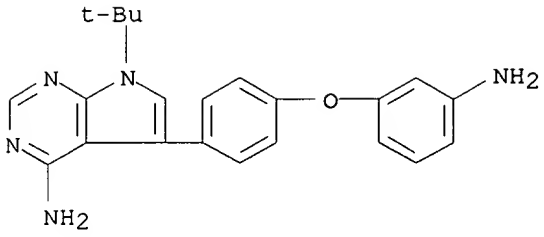
RN 213743-48-7 CAPLUS
CN Acetamide,
N-[2-[4-[4-amino-7-(1-methylethyl)-7H-pyrrolo[2,3-d]pyrimidin-5-yl]phenoxy]phenyl]- (9CI) (CA INDEX NAME)



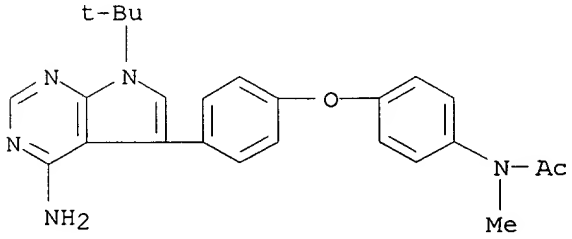
RN 213743-50-1 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 5-[4-(4-aminophenoxy)phenyl]-7-(1,1-dimethylethyl)- (9CI) (CA INDEX NAME)



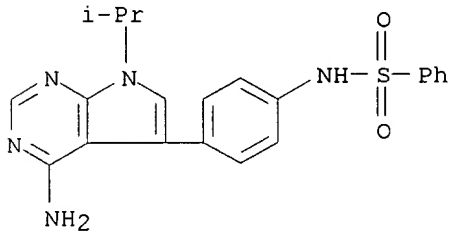
RN 213743-54-5 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 5-[4-(3-aminophenoxy)phenyl]-7-(1,1-dimethylethyl)- (9CI) (CA INDEX NAME)



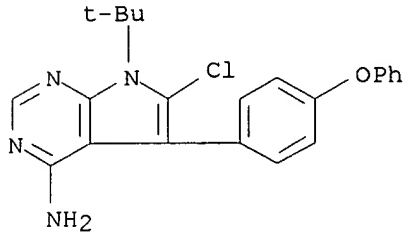
RN 213743-56-7 CAPLUS
CN Acetamide, N-[4-[4-[4-amino-7-(1,1-dimethylethyl)-7H-pyrrolo[2,3-d]pyrimidin-5-yl]phenoxy]phenyl]-N-methyl- (9CI) (CA INDEX NAME)



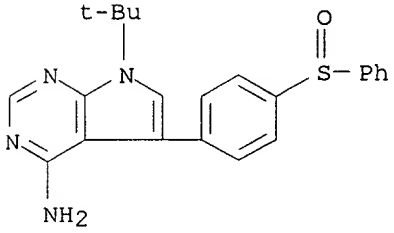
RN 213743-58-9 CAPLUS
CN Benzenesulfonamide, N-[4-[4-amino-7-(1-methylethyl)-7H-pyrrolo[2,3-d]pyrimidin-5-yl]phenyl]- (9CI) (CA INDEX NAME)



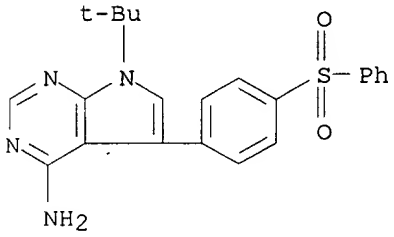
RN 213743-60-3 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 6-chloro-7-(1,1-dimethylethyl)-5-(4-phenoxyphenyl)- (9CI) (CA INDEX NAME)



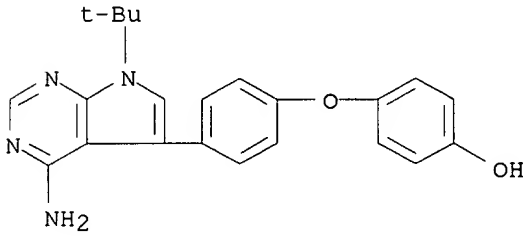
RN 213743-62-5 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 7-(1,1-dimethylethyl)-5-[4-(phenylsulfinyl)phenyl]- (9CI) (CA INDEX NAME)



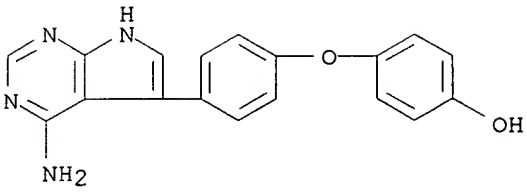
RN 213743-64-7 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 7-(1,1-dimethylethyl)-5-[4-(phenylsulfonyl)phenyl]- (9CI) (CA INDEX NAME)



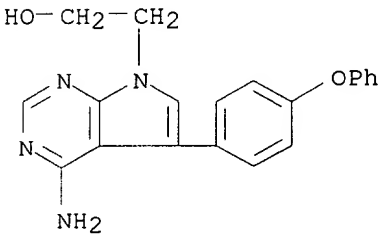
RN 213743-66-9 CAPLUS
CN Phenol, 4-[4-[4-amino-7-(1,1-dimethylethyl)-7H-pyrrolo[2,3-d]pyrimidin-5-yl]phenoxy]- (9CI) (CA INDEX NAME)



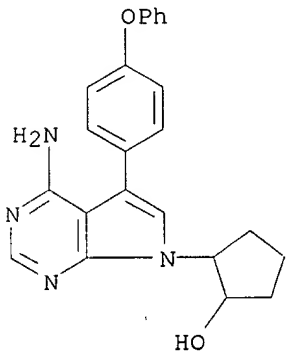
RN 213743-68-1 CAPLUS
CN Phenol, 4-[4-(4-amino-1H-pyrrolo[2,3-d]pyrimidin-5-yl)phenoxy]- (9CI) (CA
INDEX NAME)



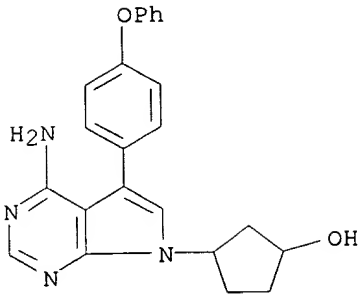
RN 213743-70-5 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidine-7-ethanol, 4-amino-5-(4-phenoxyphenyl)- (9CI) (CA INDEX NAME)



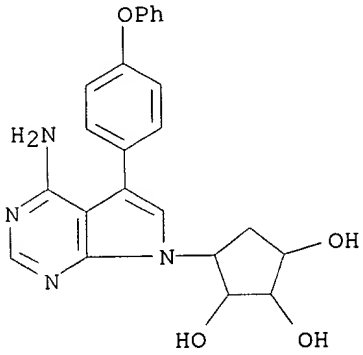
RN 213743-72-7 CAPLUS
CN Cyclopentanol, 2-[4-amino-5-(4-phenoxyphenyl)-7H-pyrrolo[2,3-d]pyrimidin-7-yl]- (9CI) (CA INDEX NAME)



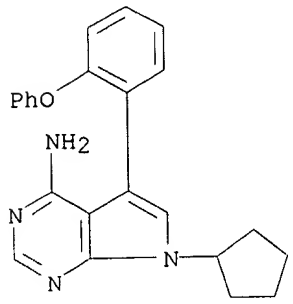
RN 213743-76-1 CAPLUS
CN Cyclopentanol, 3-[4-amino-5-(4-phenoxyphenyl)-7H-pyrrolo[2,3-d]pyrimidin-7-yl]- (9CI) (CA INDEX NAME)



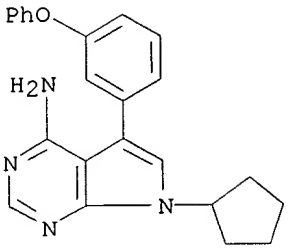
RN 213743-78-3 CAPLUS
CN 1,2,3-Cyclopentanetriol, 4-[4-amino-5-(4-phenoxyphenyl)-7H-pyrrolo[2,3-d]pyrimidin-7-yl]- (9CI) (CA INDEX NAME)



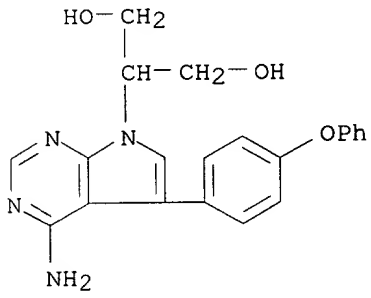
RN 213743-80-7 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 7-cyclopentyl-5-(2-phenoxyphenyl)- (9CI) (CA INDEX NAME)



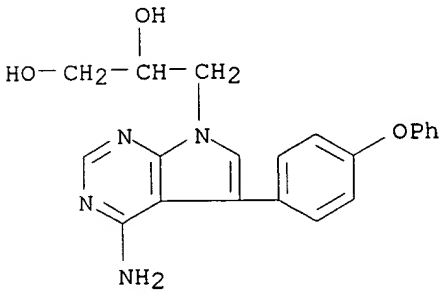
RN 213743-82-9 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 7-cyclopentyl-5-(3-phenoxyphenyl)- (9CI) (CA INDEX NAME)



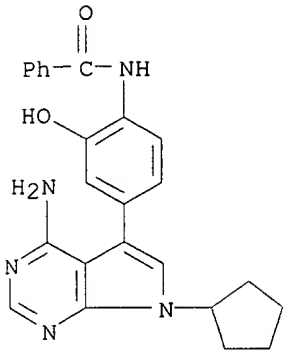
RN 213743-86-3 CAPLUS
CN 1,3-Propanediol,
2-[4-amino-5-(4-phenoxyphenyl)-7H-pyrrolo[2,3-d]pyrimidin-
7-yl]- (9CI) (CA INDEX NAME)



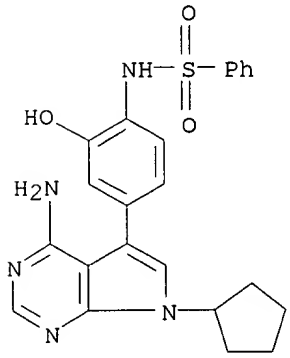
RN 213743-88-5 CAPLUS
CN 1,2-Propanediol,
3-[4-amino-5-(4-phenoxyphenyl)-7H-pyrrolo[2,3-d]pyrimidin-
7-yl]- (9CI) (CA INDEX NAME)



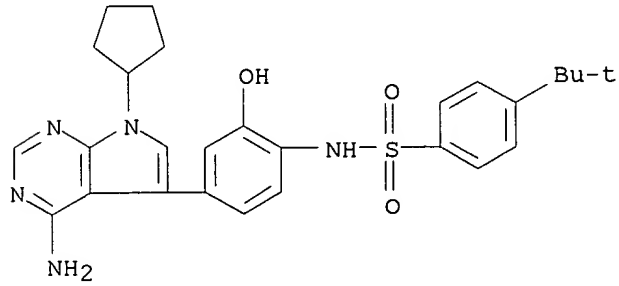
RN 213743-92-1 CAPLUS
CN Benzamide,
N-[4-(4-amino-7-cyclopentyl-7H-pyrrolo[2,3-d]pyrimidin-5-yl)-2-
hydroxyphenyl]- (9CI) (CA INDEX NAME)



RN 213743-96-5 CAPLUS
CN Benzenesulfonamide,
N-[4-(4-amino-7-cyclopentyl-7H-pyrrolo[2,3-d]pyrimidin-
5-yl)-2-hydroxyphenyl]- (9CI) (CA INDEX NAME)

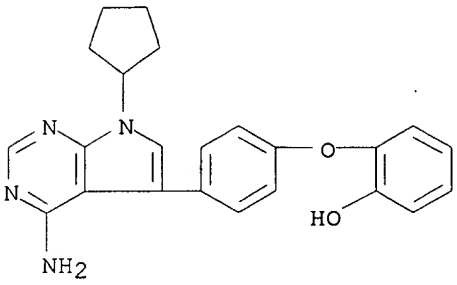


RN 213743-98-7 CAPLUS
CN Benzenesulfonamide,
N-[4-(4-amino-7-cyclopentyl-7H-pyrrolo[2,3-d]pyrimidin-
5-yl)-2-hydroxyphenyl]-4-(1,1-dimethylethyl)- (9CI) (CA INDEX NAME)

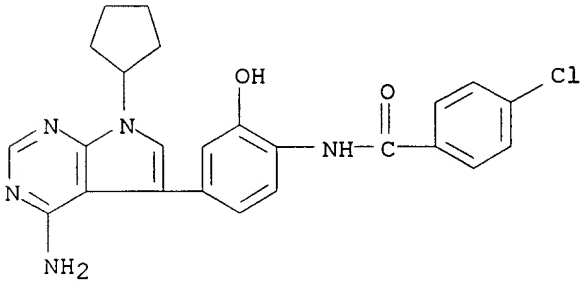


RN 213744-02-6 CAPLUS

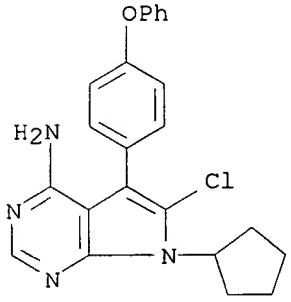
CN Phenol, 2-[4-(4-amino-7-cyclopentyl-7H-pyrrolo[2,3-d]pyrimidin-5-yl)phenoxy]- (9CI) (CA INDEX NAME)



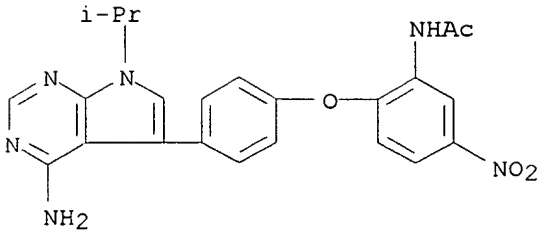
RN 213744-04-8 CAPLUS
CN Benzamide,
N-[4-(4-amino-7-cyclopentyl-7H-pyrrolo[2,3-d]pyrimidin-5-yl)-2-hydroxyphenyl]-4-chloro- (9CI) (CA INDEX NAME)



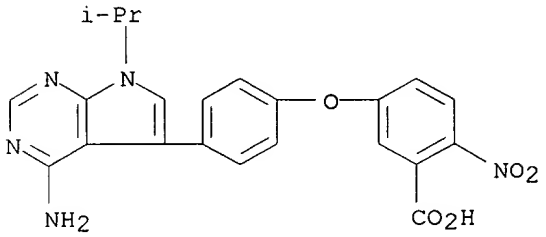
RN 213744-06-0 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 6-chloro-7-cyclopentyl-5-(4-phenoxyphenyl)- (9CI) (CA INDEX NAME)



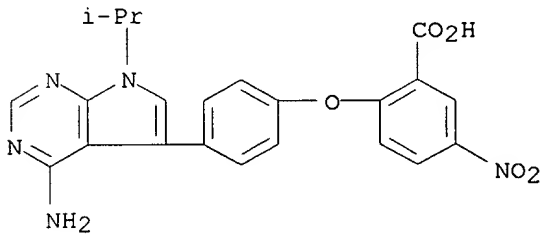
RN 213744-12-8 CAPLUS
CN Acetamide,
N-[2-[4-[4-amino-7-(1-methylethyl)-7H-pyrrolo[2,3-d]pyrimidin-5-yl]phenoxy]-5-nitrophenyl]- (9CI) (CA INDEX NAME)



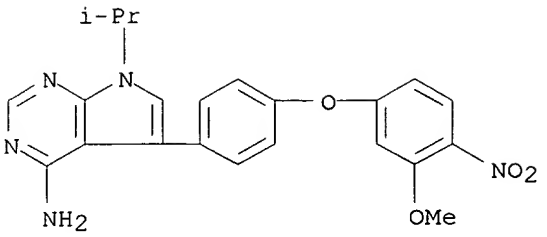
RN 213744-14-0 CAPLUS
CN Benzoic acid,
5-[4-[4-amino-7-(1-methylethyl)-7H-pyrrolo[2,3-d]pyrimidin-5-yl]phenoxy]-2-nitro- (9CI) (CA INDEX NAME)



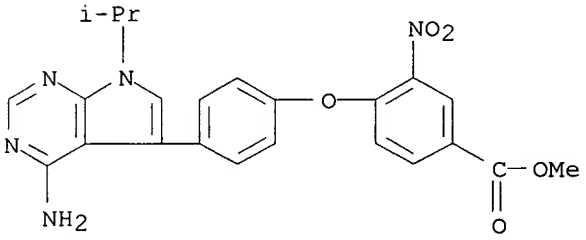
RN 213744-16-2 CAPLUS
CN Benzoic acid,
2-[4-[4-amino-7-(1-methylethyl)-7H-pyrrolo[2,3-d]pyrimidin-5-yl]phenoxy]-5-nitro- (9CI) (CA INDEX NAME)



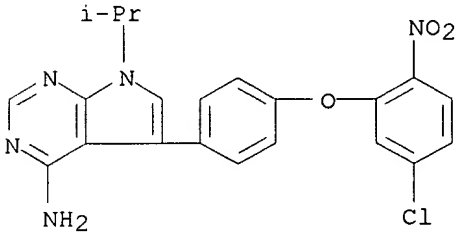
RN 213744-18-4 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine,
5-[4-(3-methoxy-4-nitrophenoxy)phenyl]-
7-(1-methylethyl)- (9CI) (CA INDEX NAME)



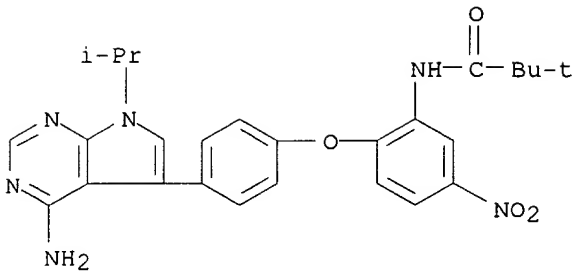
RN 213744-20-8 CAPLUS
CN Benzoic acid,
4-[4-[4-amino-7-(1-methylethyl)-7H-pyrrolo[2,3-d]pyrimidin-5-yl]phenoxy]-3-nitro-, methyl ester (9CI) (CA INDEX NAME)



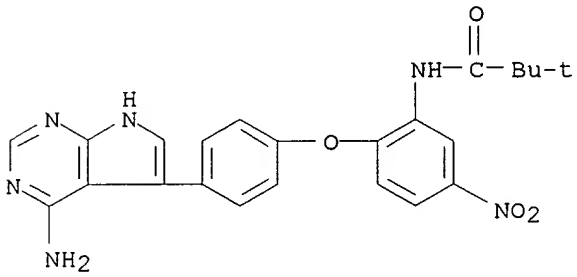
RN 213744-22-0 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine,
5-[4-(5-chloro-2-nitrophenoxy)phenyl]-7-(1-methylethyl)- (9CI) (CA INDEX NAME)



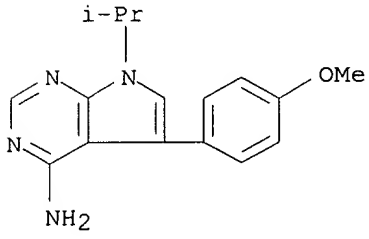
RN 213744-24-2 CAPLUS
CN Propanamide,
N-[2-[4-[4-amino-7-(1-methylethyl)-7H-pyrrolo[2,3-d]pyrimidin-5-yl]phenoxy]-5-nitrophenyl]-2,2-dimethyl- (9CI) (CA INDEX NAME)



RN 213744-29-7 CAPLUS
CN Propanamide, N-[2-[4-(4-amino-1H-pyrrolo[2,3-d]pyrimidin-5-yl)phenoxy]-5-nitrophenyl]-2,2-dimethyl- (9CI) (CA INDEX NAME)



IT **213744-87-7P**
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation)
(prepn. of pyrrolo[2,3-d]pyrimidines as tyrosine kinase inhibitors)
RN 213744-87-7 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine,
5-(4-methoxyphenyl)-7-(1-methylethyl)-
(9CI) (CA INDEX NAME)

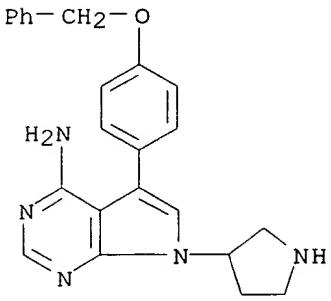


L6 ANSWER 5 OF 10 CAPLUS COPYRIGHT 2000 ACS
AN 1997:640667 CAPLUS
DN 127:318974
TI Preparation of 7-heterocyclylpyrrolo[2,3-d]pyrimidines and analogs as protein tyrosine kinase pp60c-src inhibitors
IN Altmann, Eva
PA Novartis A.-G., Switz.; Altmann, Eva
SO PCT Int. Appl., 66 pp.
CODEN: PIXXD2
DT Patent
LA German
FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	WO 9734895	A1	19970925	WO 1997-EP1095	19970305
	W: AL, AU, BA, BB, BG, BR, CA, CN, CU, CZ, EE, GE, HU, IL, IS, JP, KP, KR, LC, LK, LR, LT, LV, MG, MK, MN, MX, NO, NZ, PL, RO, SG, SI, SK, TR, TT, UA, US, UZ, VN, YU, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM				
	RW: GH, KE, LS, MW, SD, SZ, UG, AT, BE, CH, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN, ML, MR, NE, SN, TD, TG				
	CA 2249739	AA	19970925	CA 1997-2249739	19970305
	AU 9721534	A1	19971010	AU 1997-21534	19970305
	EP 888353	A1	19990107	EP 1997-914189	19970305
	R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, PT, IE, FI				
	CN 1216544	A	19990512	CN 1997-193839	19970305
	BR 9709443	A	19990810	BR 1997-9443	19970305
	NO 9804199	A	19981105	NO 1998-4199	19980911
PRAI	CH 1996-694	19960315			
	WO 1997-EP1095	19970305			

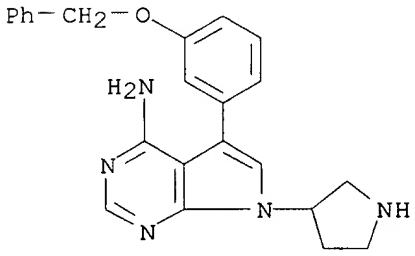
OS MARPAT 127:318974
AB Title compds. [I; R = R5Z(CH2)0-4; R1 = aryl; R2,R3 = H, halo, alkyl; R5 = H, alkyl, alkanoyl, alkoxycarbonyl, etc.; Z = (un)substituted pyrrolidine-1,2- or 1,3-diyl, -piperidine-1,2-, -1,3-, or -1,4-diyl] were prepd. as protein tyrosine kinase pp60c-src inhibitors (no data). Thus, PhCOCH2NHAc was cyclocondensed with CH2(CN)2 and the product condensed with HC(OEt)3 and NH3 to give N-(3-cyano-4-phenyl-2-pyrrolyl)formamidine which was cyclized to give, after deprotection, I (R1 = Ph, R2 = R3 = H)(II; R = H) which was condensed with Me (2R,4R)-1-tert-butoxycarbonyl-4-tosyloxypyrrolidine-2-carboxylate to give, after deprotection, II [R = (2R,4S)-2-ethoxycarbonyl-4-pyrrolidinyl].
IT **197525-41-0P 197525-42-1P 197525-44-3P 197525-45-4P 197525-49-8P 197525-50-1P 197525-53-4P 197525-54-5P 197525-55-6P 197525-56-7P 197525-57-8P**
RL: BAC (Biological activity or effector, except adverse); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
(prepn. of 7-heterocyclylpyrrolo[2,3-d]pyrimidines and analogs as protein tyrosine kinase pp60c-src inhibitors)
RN 197525-41-0 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 5-[4-(phenylmethoxy)phenyl]-7-(3-

pyrrolidinyl)-, dihydrochloride (9CI) (CA INDEX NAME)



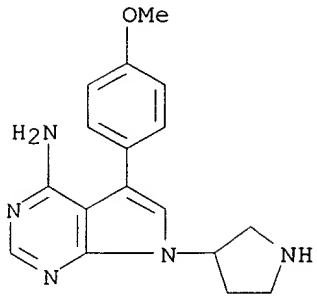
● 2 HCl

RN 197525-42-1 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 5-[3-(phenylmethoxy)phenyl]-7-(3-pyrrolidinyl)-, dihydrochloride (9CI) (CA INDEX NAME)



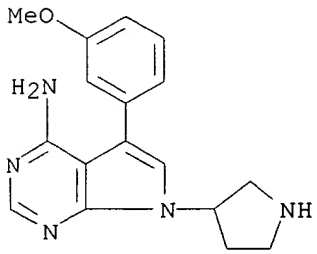
● 2 HCl

RN 197525-44-3 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 5-(4-methoxyphenyl)-7-(3-pyrrolidinyl)-, dihydrochloride (9CI) (CA INDEX NAME)



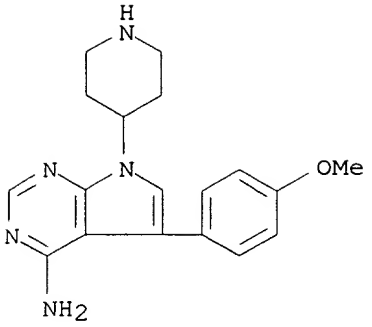
●2 HCl

RN 197525-45-4 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 5-(3-methoxyphenyl)-7-(3-pyrrolidinyl)-, dihydrochloride (9CI) (CA INDEX NAME)

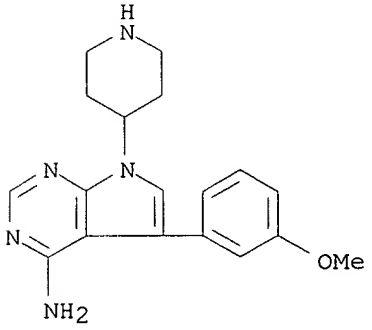


●2 HCl

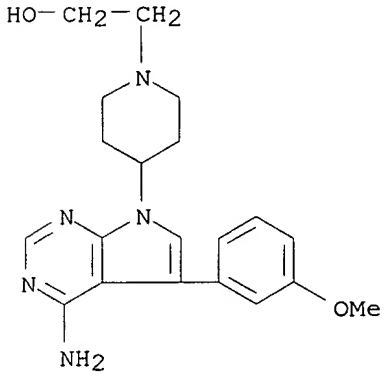
RN 197525-49-8 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 5-(4-methoxyphenyl)-7-(4-piperidinyl)-(9CI) (CA INDEX NAME)



RN 197525-50-1 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 5-(3-methoxyphenyl)-7-(4-piperidinyl)- (9CI) (CA INDEX NAME)

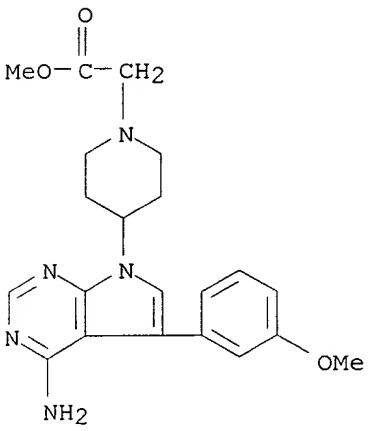


RN 197525-53-4 CAPLUS
CN 1-Piperidineethanol, 4-[4-amino-5-(3-methoxyphenyl)-7H-pyrrolo[2,3-d]pyrimidin-7-yl]- (9CI) (CA INDEX NAME)

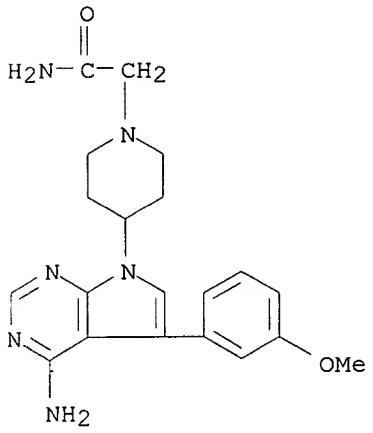


RN 197525-54-5 CAPLUS

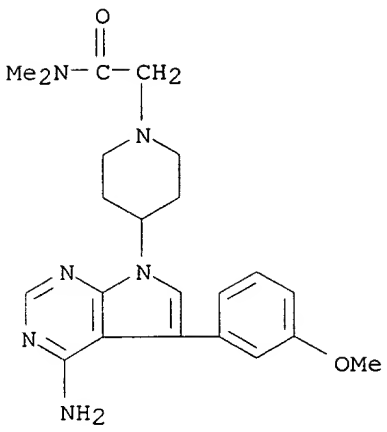
CN 1-Piperidineacetic acid, 4-[4-amino-5-(3-methoxyphenyl)-7H-pyrrolo[2,3-d]pyrimidin-7-yl]-, methyl ester (9CI) (CA INDEX NAME)



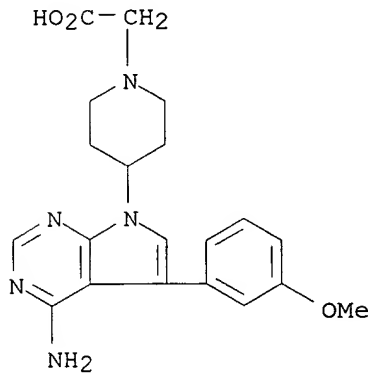
RN 197525-55-6 CAPLUS
CN 1-Piperidineacetamide, 4-[4-amino-5-(3-methoxyphenyl)-7H-pyrrolo[2,3-d]pyrimidin-7-yl]- (9CI) (CA INDEX NAME)



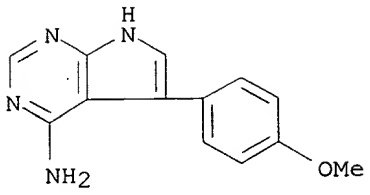
RN 197525-56-7 CAPLUS
CN 1-Piperidineacetamide, 4-[4-amino-5-(3-methoxyphenyl)-7H-pyrrolo[2,3-d]pyrimidin-7-yl]-N,N-dimethyl- (9CI) (CA INDEX NAME)



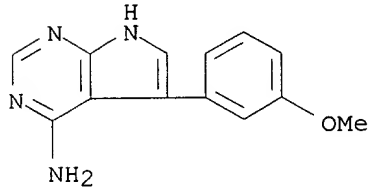
RN 197525-57-8 CAPLUS
CN 1-Piperidineacetic acid, 4-[4-amino-5-(3-methoxyphenyl)-7H-pyrrolo[2,3-d]pyrimidin-7-yl]- (9CI) (CA INDEX NAME)



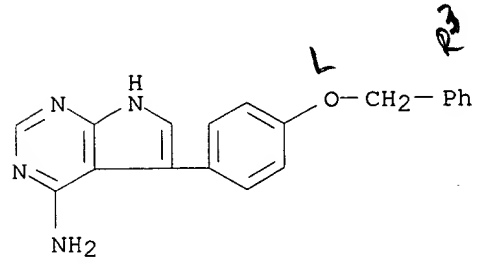
IT **121405-37-6P 194787-34-3P 194787-35-4P**
194787-36-5P 197525-66-9P 197525-67-0P
197525-68-1P 197525-83-0P 197525-84-1P
197525-86-3P 197525-87-4P
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation)
(prepn. of 7-heterocyclylpyrrolo[2,3-d]pyrimidines and analogs as protein tyrosine kinase pp60c-src inhibitors)
RN 121405-37-6 CAPLUS
CN 1H-Pyrrolo[2,3-d]pyrimidin-4-amine, 5-(4-methoxyphenyl)- (9CI) (CA INDEX NAME)



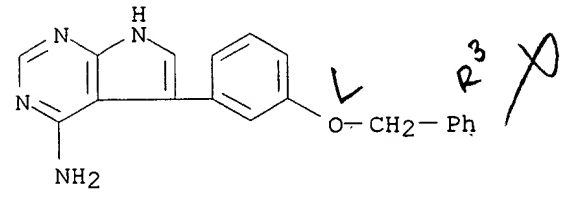
RN 194787-34-3 CAPLUS
CN 1H-Pyrrolo[2,3-d]pyrimidin-4-amine, 5-(3-methoxyphenyl)- (9CI) (CA INDEX NAME)



RN 194787-35-4 CAPLUS
CN 1H-Pyrrolo[2,3-d]pyrimidin-4-amine, 5-[4-(phenylmethoxy)phenyl]- (9CI) (CA INDEX NAME)

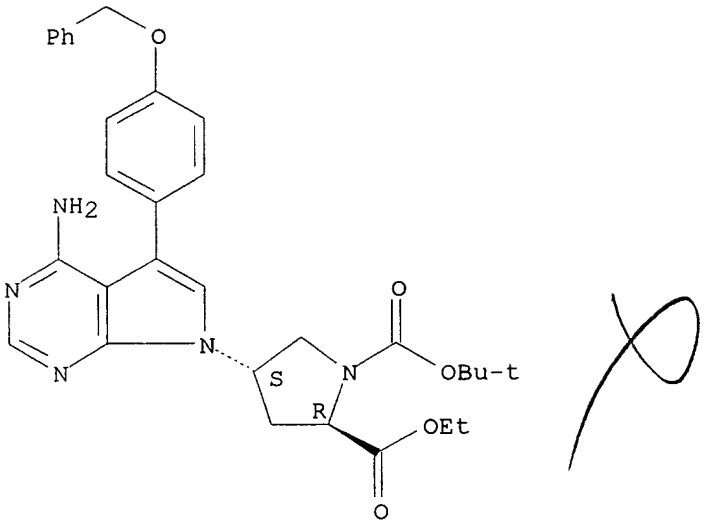


RN 194787-36-5 CAPLUS
CN 1H-Pyrrolo[2,3-d]pyrimidin-4-amine, 5-[3-(phenylmethoxy)phenyl]- (9CI) (CA INDEX NAME)



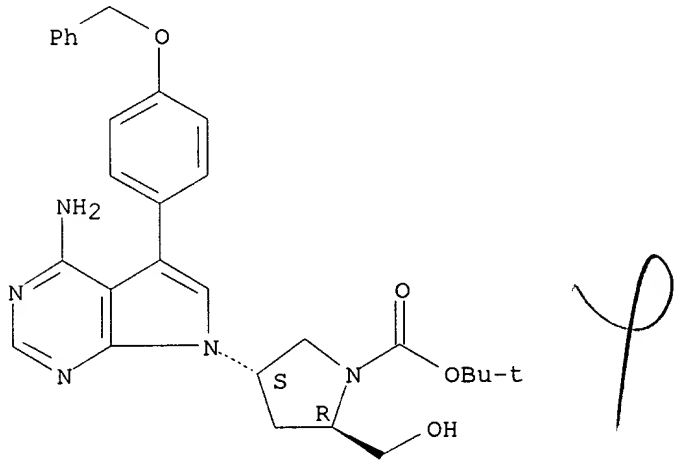
RN 197525-66-9 CAPLUS
CN 1,2-Pyrrolidinedicarboxylic acid, 4-[4-amino-5-[4-(phenylmethoxy)phenyl]-7H-pyrrolo[2,3-d]pyrimidin-7-yl]-, 1-(1,1-dimethylethyl) 2-ethyl ester, (2R-trans)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



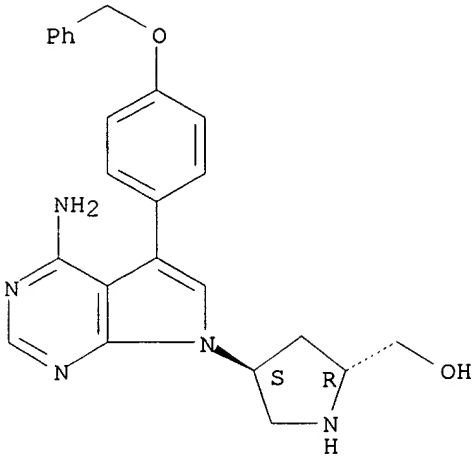
RN 197525-67-0 CAPLUS
CN 1-Pyrrolidinecarboxylic acid, 4-[4-amino-5-[4-(phenylmethoxy)phenyl]-7H-pyrrolo[2,3-d]pyrimidin-7-yl]-2-(hydroxymethyl)-, 1,1-dimethylethyl ester,
(2R-trans)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

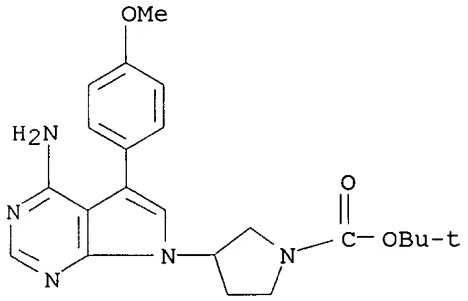


RN 197525-68-1 CAPLUS
CN 2-Pyrrolidinemethanol, 4-[4-amino-5-[4-(phenylmethoxy)phenyl]-7H-pyrrolo[2,3-d]pyrimidin-7-yl]-, (2R-trans)- (9CI) (CA INDEX NAME)

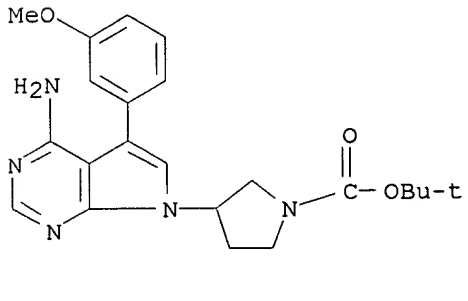
Absolute stereochemistry.



RN 197525-83-0 CAPLUS
CN 1-Pyrrolidinecarboxylic acid, 3-[4-amino-5-(4-methoxyphenyl)-7H-pyrrolo[2,3-d]pyrimidin-7-yl]-, 1,1-dimethylethyl ester (9CI) (CA INDEX NAME)

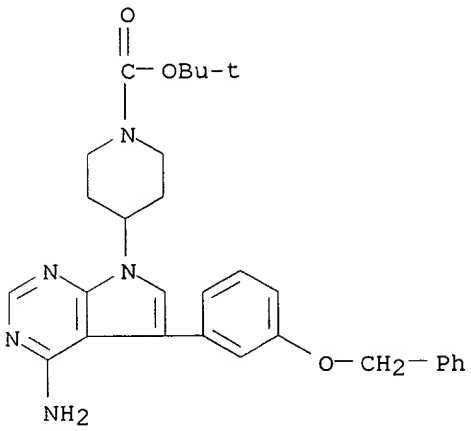


RN 197525-84-1 CAPLUS
CN 1-Pyrrolidinecarboxylic acid, 3-[4-amino-5-(3-methoxyphenyl)-7H-pyrrolo[2,3-d]pyrimidin-7-yl]-, 1,1-dimethylethyl ester (9CI) (CA INDEX NAME)

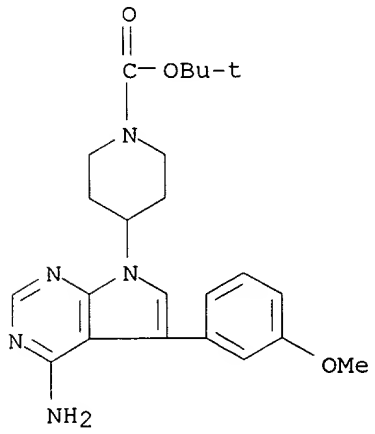


RN 197525-86-3 CAPLUS
CN 1-Piperidinecarboxylic acid, 4-[4-amino-5-[3-(phenylmethoxy)phenyl]-7H-

pyrrolo[2,3-d]pyrimidin-7-yl]-, 1,1-dimethylethyl ester (9CI) (CA INDEX NAME)



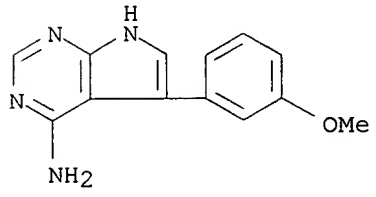
RN 197525-87-4 CAPLUS
CN 1-Piperidinecarboxylic acid, 4-[4-amino-5-(3-methoxyphenyl)-7H-pyrrolo[2,3-d]pyrimidin-7-yl]-, 1,1-dimethylethyl ester (9CI) (CA INDEX NAME)



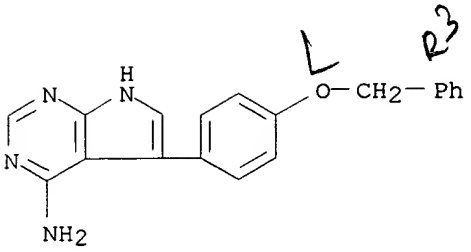
L6 ANSWER 6 OF 10 CAPLUS COPYRIGHT 2000 ACS
AN 1997:618099 CAPLUS
DN 127:293238
TI 7-Alkylpyrrolo[2,3-d]pyrimidines
IN Altmann, Eva; Missbach, Martin; Widler, Leo
PA Novartis A.-G., Switz.; Altmann, Eva; Missbach, Martin; Widler, Leo
SO PCT Int. Appl., 64 pp.
CODEN: PIXXD2
DT Patent
LA German
FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	WO 9732879	A1	19970912	WO 1997-EP874	19970224
	W: AL, AU, BA, BB, BG, BR, CA, CN, CU, CZ, EE, GE, HU, IL, IS, JP, KP, KR, LC, LK, LR, LT, LV, MG, MK, MN, MX, NO, NZ, PL, RO, SG, SI, SK, TR, TT, UA, US, UZ, VN, YU, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM				
	RW: KE, LS, MW, SD, SZ, UG, AT, BE, CH, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN, ML, MR, NE, SN, TD, TG				
	AU 9717946	A1	19970922	AU 1997-17946	19970224

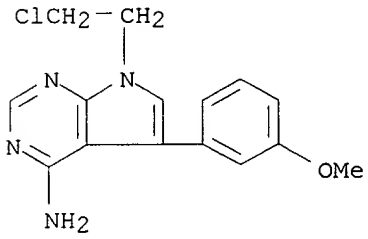
PRAI EP 1996-810137 19960306
WO 1997-EP874 19970224
OS MARPAT 127:293238
AB Approx. 50 protein tyrosine kinase inhibiting title compds. I [R = H, alkyl, hydroxyalkyl, hydroxy(aryl)alkyl, hydroxycycloalkyl, hydroxycycloalkylalkyl, chloroalkyl, substituted-aminoalkyl, etc.; R1 = Ph, substituted-phenyl] were prepd. E.g., I (R = H, R1 = Ph) was prepd. from phenacylamine-HCl and Ac2O in 4 steps via 2-amino-4-phenyl-1H-pyrrole-3-carbonitrile and N-(3-cyano-4-phenyl-1H-pyrrol-2-yl)formamidine.
IT **194787-34-3P 194787-35-4P 196964-56-4P 196964-57-5P**
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation)
(prepn. of protein tyrosine kinase inhibiting pyrrolopyrimidines)
RN 194787-34-3 CAPLUS
CN 1H-Pyrrolo[2,3-d]pyrimidin-4-amine, 5-(3-methoxyphenyl)- (9CI) (CA INDEX NAME)



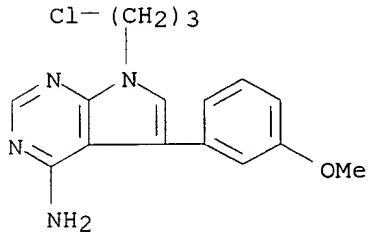
RN 194787-35-4 CAPLUS
CN 1H-Pyrrolo[2,3-d]pyrimidin-4-amine, 5-[4-(phenylmethoxy)phenyl]- (9CI) (CA INDEX NAME)



RN 196964-56-4 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 7-(2-chloroethyl)-5-(3-methoxyphenyl)- (9CI) (CA INDEX NAME)

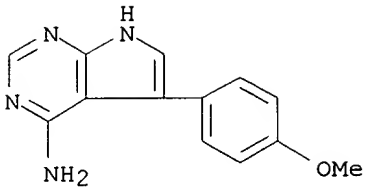


RN 196964-57-5 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 7-(3-chloropropyl)-5-(3-methoxyphenyl)- (9CI) (CA INDEX NAME)

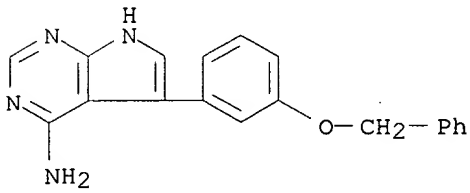


IT **121405-37-6P 194787-36-5P 196963-91-4P**
196964-10-0P 196964-14-4P 196964-19-9P
196964-28-0P 196964-29-1P 196964-30-4P
196964-31-5P 196964-32-6P 196964-64-4P
RL: SPN (Synthetic preparation); PREP (Preparation)
(prepn. of protein tyrosine kinase inhibiting pyrrolopyrimidines)

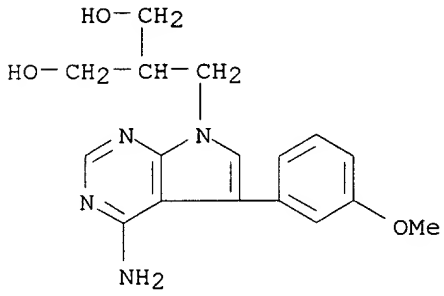
RN 121405-37-6 CAPLUS
CN 1H-Pyrrolo[2,3-d]pyrimidin-4-amine, 5-(4-methoxyphenyl)- (9CI) (CA INDEX NAME)



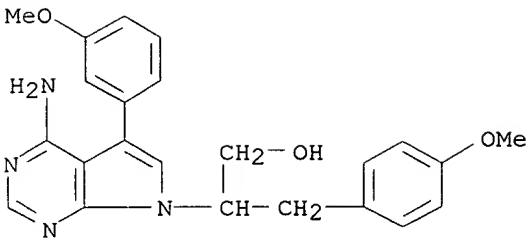
RN 194787-36-5 CAPLUS
CN 1H-Pyrrolo[2,3-d]pyrimidin-4-amine, 5-[3-(phenylmethoxy)phenyl]- (9CI) (CA INDEX NAME)



RN 196963-91-4 CAPLUS
CN 1,3-Propanediol, 2-[[4-amino-5-(3-methoxyphenyl)-7H-pyrrolo[2,3-d]pyrimidin-7-yl]methyl]- (9CI) (CA INDEX NAME)

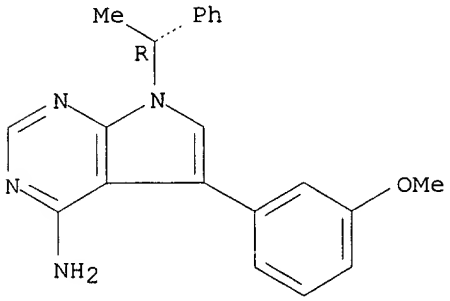


RN 196964-10-0 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidine-7-ethanol, 4-amino-5-(3-methoxyphenyl)-.beta.-[(4-methoxyphenyl)methyl]- (9CI) (CA INDEX NAME)

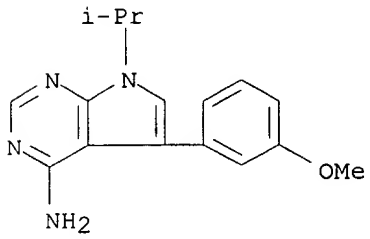


RN 196964-14-4 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 5-(3-methoxyphenyl)-7-(1-phenylethyl)-, (R)- (9CI) (CA INDEX NAME)

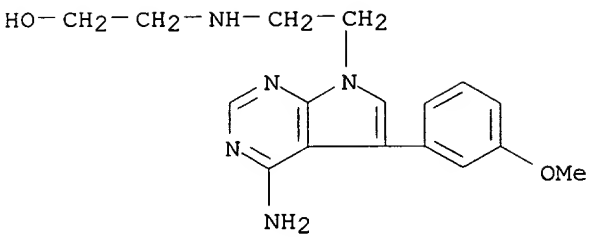
Absolute stereochemistry.



RN 196964-19-9 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 5-(3-methoxyphenyl)-7-(1-methylethyl)- (9CI) (CA INDEX NAME)

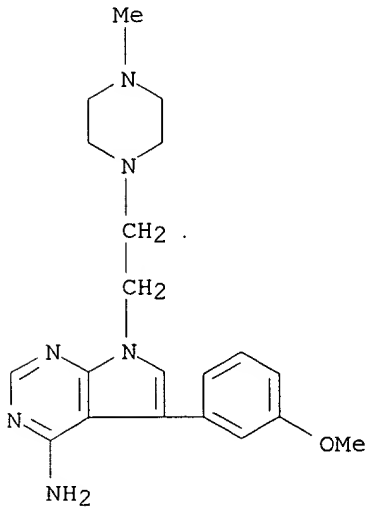


RN 196964-28-0 CAPLUS
CN Ethanol, 2-[[2-[4-amino-5-(3-methoxyphenyl)-7H-pyrrolo[2,3-d]pyrimidin-7-yl]ethyl]amino]-, dihydrochloride (9CI) (CA INDEX NAME)

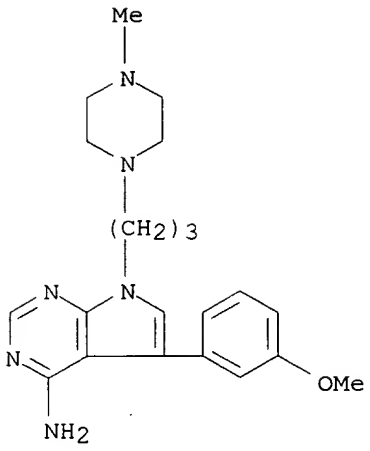


●2 HCl

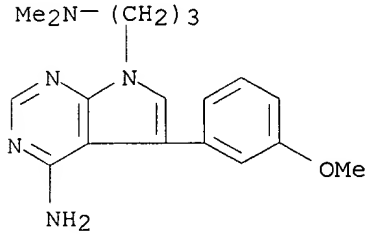
RN 196964-29-1 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 5-(3-methoxyphenyl)-7-[2-(4-methyl-1-piperazinyl)ethyl]- (9CI) (CA INDEX NAME)



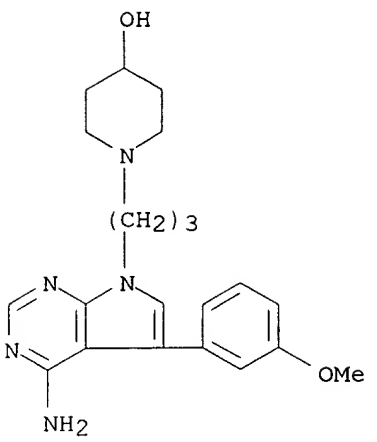
RN 196964-30-4 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 5-(3-methoxyphenyl)-7-[3-(4-methyl-1-piperazinyl)propyl]- (9CI) (CA INDEX NAME)



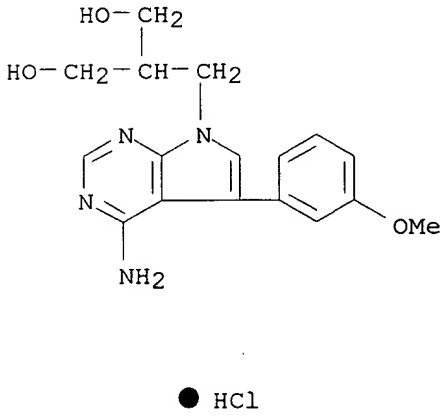
RN 196964-31-5 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidine-7-propanamine, 4-amino-5-(3-methoxyphenyl)-N,N-dimethyl- (9CI) (CA INDEX NAME)



RN 196964-32-6 CAPLUS
CN 4-Piperidinol, 1-[3-[4-amino-5-(3-methoxyphenyl)-7H-pyrrolo[2,3-d]pyrimidin-7-yl]propyl]- (9CI) (CA INDEX NAME)



RN 196964-64-4 CAPLUS
CN 1,3-Propanediol, 2-[[4-amino-5-(3-methoxyphenyl)-7H-pyrrolo[2,3-d]pyrimidin-7-yl]methyl]-, monohydrochloride (9CI) (CA INDEX NAME)

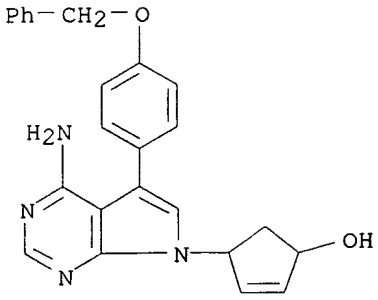


● HCl

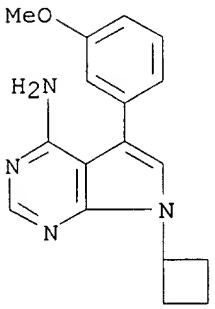
L6 ANSWER 7 OF 10 CAPLUS COPYRIGHT 2000 ACS
AN 1997:533648 CAPLUS
DN 127:220666
TI Preparation of 4-aminopyrrolo[2,3-d]pyrimidines as tyrosine kinase inhibitors
IN Altmann, Eva; Widler, Leo; Missbach, Martin
PA Novartis A.-G, Switz.; Altmann, Eva; Widler, Leo; Missbach, Martin
SO PCT Int. Appl., 48 pp.
CODEN: PIXXD2
DT Patent
LA English
FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	WO 9728161	A1	19970807	WO 1997-EP350	19970127
	W: AL, AU, BA, BB, BG, BR, CA, CN, CU, CZ, EE, GE, HU, IL, IS, JP, KP, KR, LC, LK, LR, LT, LV, MG, MK, MN, MX, NO, NZ, PL, RO, SG, SI, SK, TR, TT, UA, US, UZ, VN, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM				
	RW: KE, LS, MW, SD, SZ, UG, AT, BE, CH, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN, ML, MR, NE, SN, TD, TG				
	AU 9715951	A1	19970822	AU 1997-15951	19970127
PRAI	CH 1996-257		19960201		
	WO 1997-EP350		19970127		

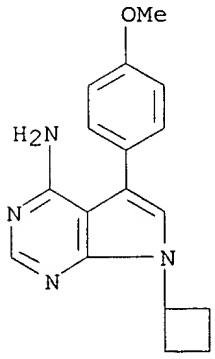
OS MARPAT 127:220666
AB Title compds. [I; R1 = aryl; R2,R4 = H, halo, alkyl; R3 = (un)substituted cycloalk(en)yl] were prepd. as tyrosine kinase inhibitors (no data). Thus, PhCOCH2NH2 was N-acetylated and the product cyclocondensed with CH2(CN)2 to give 2-amino-3-cyano-4-phenylpyrrole which was condensed with HC(OEt)3 and NH3 and the product cyclized to give I (R1 = Ph, R2 = R4 = H)(II; R3 = H). The latter was condensed with cyclopentadiene monoepoxide to give II (R3 = 3-hydroxy-2-cyclopentenyl).
IT **194787-38-7P 194787-74-1P 194787-80-9P 194787-83-2P 194787-84-3P 194787-86-5P 194787-87-6P**
RL: BAC (Biological activity or effector, except adverse); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
(prepn. of 4-aminopyrrolo[2,3-d]pyrimidines as tyrosine kinase inhibitors)
RN 194787-38-7 CAPLUS
CN 2-Cyclopenten-1-ol, 4-[4-amino-5-[4-(phenylmethoxy)phenyl]-7H-pyrrolo[2,3-d]pyrimidin-7-yl]- (9CI) (CA INDEX NAME)



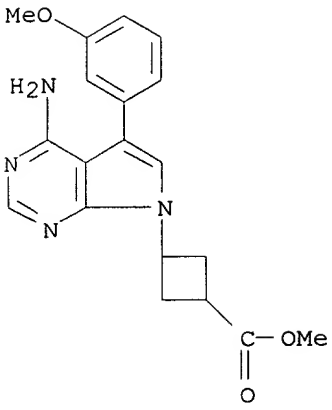
RN 194787-74-1 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 7-cyclobutyl-5-(3-methoxyphenyl)-(9CI) (CA INDEX NAME)



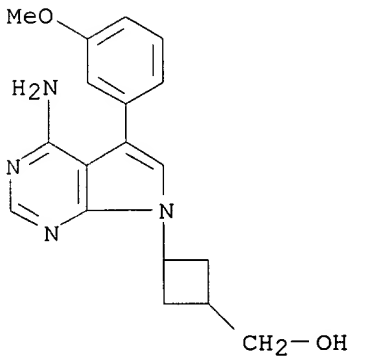
RN 194787-80-9 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 7-cyclobutyl-5-(4-methoxyphenyl)-(9CI) (CA INDEX NAME)



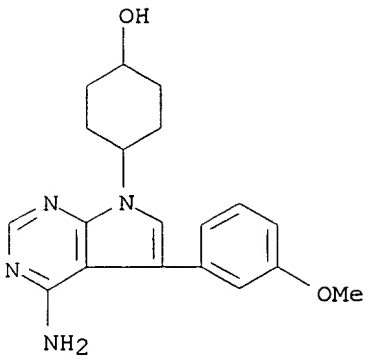
RN 194787-83-2 CAPLUS
CN Cyclobutanecarboxylic acid, 3-[4-amino-5-(3-methoxyphenyl)-7H-pyrrolo[2,3-d]pyrimidin-7-yl]-, methyl ester (9CI) (CA INDEX NAME)



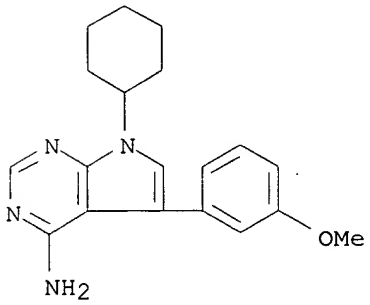
RN 194787-84-3 CAPLUS
CN Cyclobutanemethanol, 3-[4-amino-5-(3-methoxyphenyl)-7H-pyrrolo[2,3-d]pyrimidin-7-yl]- (9CI) (CA INDEX NAME)



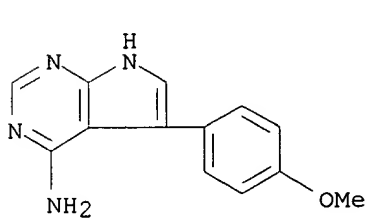
RN 194787-86-5 CAPLUS
CN Cyclohexanol, 4-[4-amino-5-(3-methoxyphenyl)-7H-pyrrolo[2,3-d]pyrimidin-7-yl]- (9CI) (CA INDEX NAME)



RN 194787-87-6 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 7-cyclohexyl-5-(3-methoxyphenyl)- (9CI) (CA INDEX NAME)

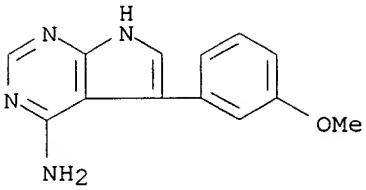


IT **121405-37-6P 194787-34-3P 194787-35-4P**
194787-36-5P 194788-04-0P 194788-05-1P
194788-06-2P
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation)
(prepn. of 4-aminopyrrolo[2,3-d]pyrimidines as tyrosine kinase inhibitors)
RN 121405-37-6 CAPLUS
CN 1H-Pyrrolo[2,3-d]pyrimidin-4-amine, 5-(4-methoxyphenyl)- (9CI) (CA INDEX NAME)

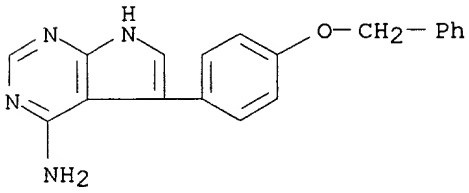


RN 194787-34-3 CAPLUS
CN 1H-Pyrrolo[2,3-d]pyrimidin-4-amine, 5-(3-methoxyphenyl)- (9CI) (CA INDEX

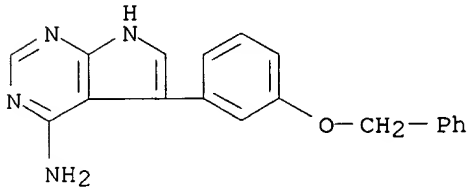
NAME)



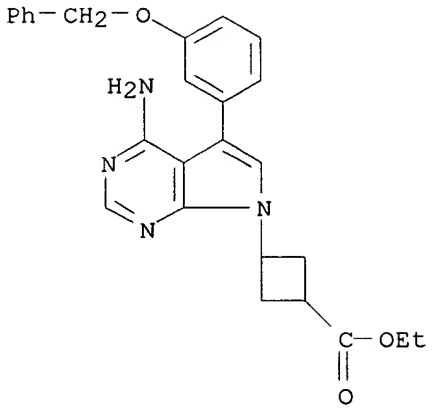
RN 194787-35-4 CAPLUS
CN 1H-Pyrrolo[2,3-d]pyrimidin-4-amine, 5-[4-(phenylmethoxy)phenyl]- (9CI) (CA INDEX NAME)



RN 194787-36-5 CAPLUS
CN 1H-Pyrrolo[2,3-d]pyrimidin-4-amine, 5-[3-(phenylmethoxy)phenyl]- (9CI) (CA INDEX NAME)

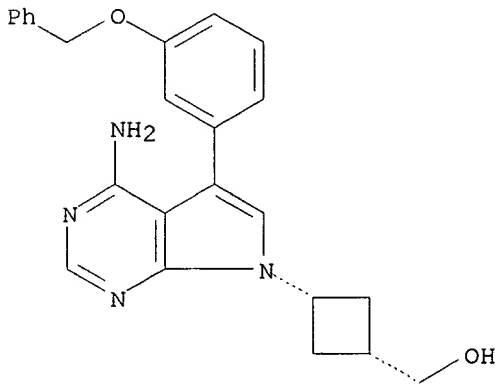


RN 194788-04-0 CAPLUS
CN Cyclobutanecarboxylic acid, 3-[4-amino-5-[3-(phenylmethoxy)phenyl]-7H-pyrrolo[2,3-d]pyrimidin-7-yl]-, ethyl ester (9CI) (CA INDEX NAME)



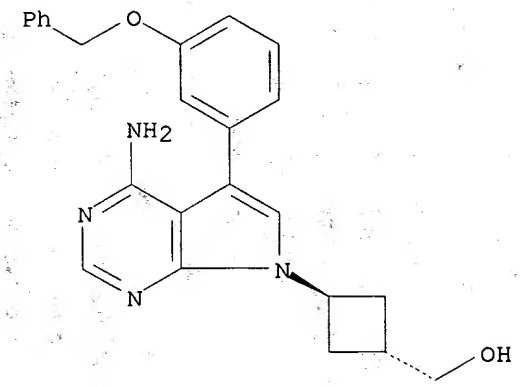
RN 194788-05-1 CAPLUS
CN Cyclobutanemethanol,
3-[4-amino-5-[3-(phenylmethoxy)phenyl]-7H-pyrrolo[2,3-
d]pyrimidin-7-yl]-, cis- (9CI) (CA INDEX NAME)

Relative stereochemistry.



RN 194788-06-2 CAPLUS
CN Cyclobutanemethanol,
3-[4-amino-5-[3-(phenylmethoxy)phenyl]-7H-pyrrolo[2,3-
d]pyrimidin-7-yl]-, trans- (9CI) (CA INDEX NAME)

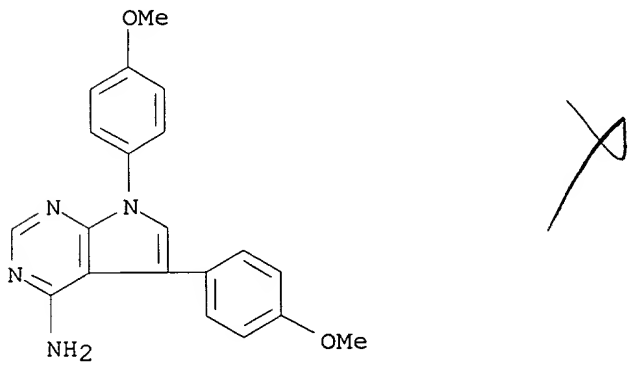
Relative stereochemistry.


Ph
O
NH2
N
N
N
OH

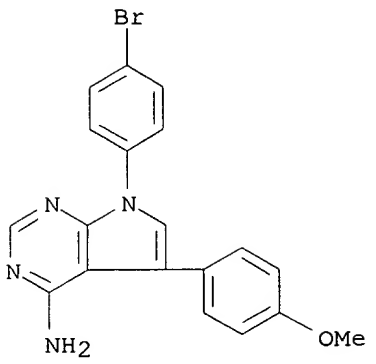
L6 ANSWER 8 OF 10 CAPLUS COPYRIGHT 2000 ACS
AN 1997:411066 CAPLUS
DN 127:90501
TI 4-aminopyrrolo[2,3-d]pyrimidines as tyrosine kinase inhibitors
IN Dow, Robert L.; Koch, Kevin
PA Pfizer Inc., USA
SO U.S., 5 pp.
CODEN: USXXAM
DT Patent
LA English
FAN.CNT 1

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
US 5639757	A	19970617	US 1995-448248	19950523

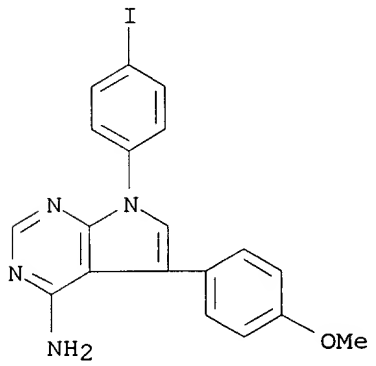
OS MARPAT 127:90501
AB Certain 4-aminopyrrolo[2,3-d]pyrimidine compds., and their pharmaceutically-acceptable salts (Markush structure given), are inhibitors of tyrosine kinase enzymes, and are useful for immunoregulation and for the treatment of cancer, angiogenesis and atherosclerosis. Inhibitory activity of 15 title compds. at 1x10-4-10-6 M concn. was tested on enzyme pp60src, a tyrosine-specific phosphokinase assocd. with the inner surface of the plasma membrane.
IT **121405-31-0 121405-32-1 121405-33-2**
RL: BAC (Biological activity or effector, except adverse); THU (Therapeutic use); BIOL (Biological study); USES (Uses)
(aminopyrrolopyrimidines as tyrosine kinase inhibitors)
RN 121405-31-0 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 5,7-bis(4-methoxyphenyl)- (9CI) (CA INDEX NAME)



RN 121405-32-1 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 7-(4-bromophenyl)-5-(4-methoxyphenyl)- (9CI) (CA INDEX NAME)



RN 121405-33-2 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 7-(4-iodophenyl)-5-(4-methoxyphenyl)-
(9CI) (CA INDEX NAME)

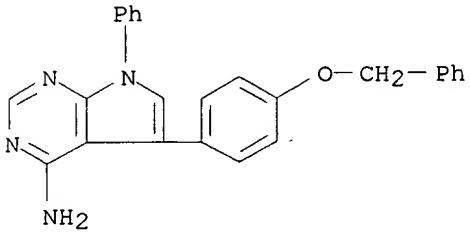


L6 ANSWER 9 OF 10 CAPLUS COPYRIGHT 2000 ACS
AN 1996:443907 CAPLUS
DN 125:86670
TI Preparation of 4-aminopyrrolo[2,3-d]pyrimidines as inhibitors of the protein tyrosine kinase pp60c-src
IN Missbach, Martin
PA Ciba-Geigy A.-G., Switz.
SO PCT Int. Appl., 77 pp.
CODEN: PIXXD2
DT Patent
LA English
FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	WO 9610028	A1	19960404	WO 1995-EP3536	19950908
	W: AM, AU, BB, BG, BR, BY, CA, CN, CZ, EE, FI, GE, HU, IS, JP, KG, KP, KR, KZ, LK, LR, LT, LV, MD, MG, MK, MN, MX, NO, NZ, PL, RO, RU, SG, SI, SK, TJ, TM, TT, UA, US, UZ, VN				
	RW: KE, MW, SD, SZ, UG, AT, BE, CH, DE, DK, ES, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN, ML, MR, NE, SN, TD, TG				
	CA 2200210	AA	19960404	CA 1995-2200210	19950908
	AU 9535643	A1	19960419	AU 1995-35643	19950908
	AU 694801	B2	19980730		
	EP 783505	A1	19970716	EP 1995-932693	19950908
	R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IE, IT, LI, LU, MC, NL, PT, SE				
	CN 1164234	A	19971105	CN 1995-196308	19950908
	CN 1046731	B	19991124		
	HU 76785	A2	19971128	HU 1997-1333	19950908
	BR 9509048	A	19980106	BR 1995-9048	19950908
	JP 10506624	T2	19980630	JP 1995-511312	19950908
	US 5869485	A	19990209	US 1997-793313	19970319
	NO 9701342	A	19970321	NO 1997-1342	19970321
	FI 9701225	A	19970514	FI 1997-1225	19970324
PRAI	CH 1994-2953	19940929			
	WO 1995-EP3536	19950908			

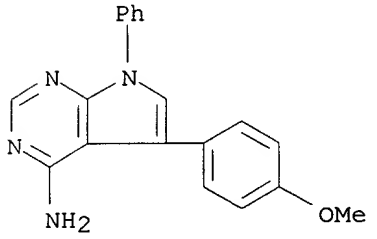
OS CASREACT 125:86670; MARPAT 125:86670
AB The title compds. [I; R1, R3 = (substituted) aryl; R2 = H, alkyl, halo], useful for the treatment of osteoporosis, breast cancer and cardiovascular disorders, e.g. thrombosis, were prepd. by e.g. treatment of substituted 2-amino-3-cyano-pyrrole with (EtO)3CH followed by treatment of II (X = EtO) with NH3/EtOH and cyclization of II (X = NH2) with NH3/EtOH at 130.degree. in an autoclave. Tablets formulations contg. I are given. In general, compds. I showed IC50 of 0.001-10 .mu.M against protein tyrosine kinase pp60c-src.
IT **178909-50-7P**
RL: BAC (Biological activity or effector, except adverse); RCT (Reactant); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
(prepn. of 4-aminopyrrolo[2,3-d]pyrimidines as inhibitors of the protein tyrosine kinase pp60c-src)
RN 178909-50-7 CAPLUS

CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 7-phenyl-5-[4-(phenylmethoxy)phenyl]- (9CI) (CA INDEX NAME)

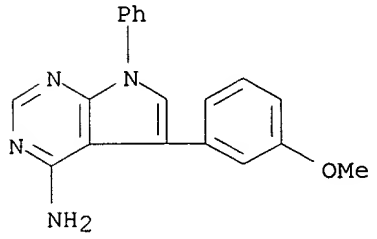


IT **121405-30-9P 178909-27-8P 178909-32-5P**
178909-42-7P 178909-46-1P 178909-47-2P
178909-49-4P 178909-66-5P 178909-84-7P
178909-97-2P 178909-98-3P 178910-00-4P
RL: BAC (Biological activity or effector, except adverse); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
(prepn. of 4-aminopyrrolo[2,3-d]pyrimidines as inhibitors of the protein tyrosine kinase pp60c-src)

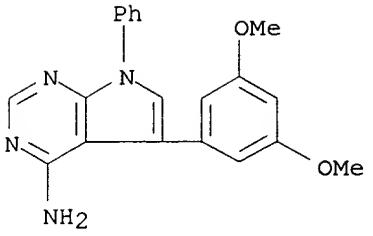
RN 121405-30-9 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 5-(4-methoxyphenyl)-7-phenyl- (9CI) (CA INDEX NAME)



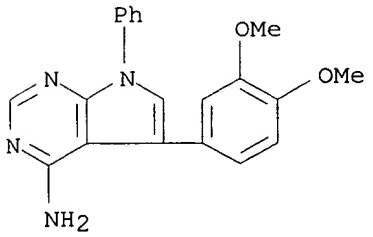
RN 178909-27-8 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 5-(3-methoxyphenyl)-7-phenyl- (9CI) (CA INDEX NAME)



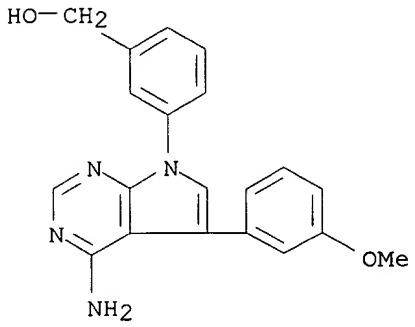
RN 178909-32-5 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 5-(3,5-dimethoxyphenyl)-7-phenyl- (9CI) (CA INDEX NAME)



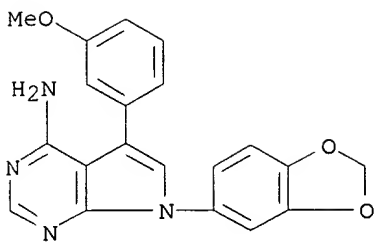
RN 178909-42-7 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 5-(3,4-dimethoxyphenyl)-7-phenyl- (9CI) (CA INDEX NAME)



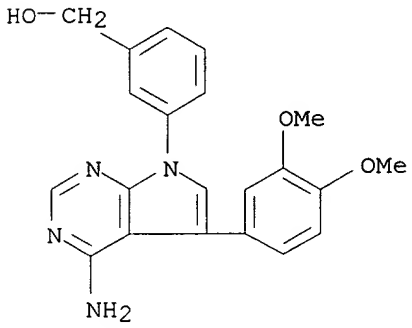
RN 178909-46-1 CAPLUS
CN Benzenemethanol, 3-[4-amino-5-(3-methoxyphenyl)-7H-pyrrolo[2,3-d]pyrimidin-7-yl]- (9CI) (CA INDEX NAME)



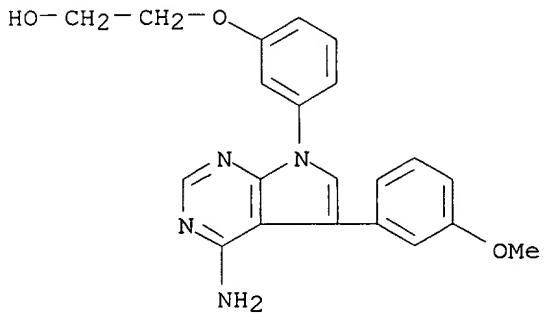
RN 178909-47-2 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 7-(1,3-benzodioxol-5-yl)-5-(3-methoxyphenyl)- (9CI) (CA INDEX NAME)



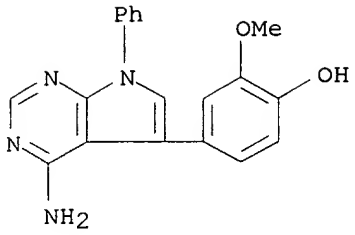
RN 178909-49-4 CAPLUS
CN Benzenemethanol, 3-[4-amino-5-(3,4-dimethoxyphenyl)-7H-pyrrolo[2,3-d]pyrimidin-7-yl]- (9CI) (CA INDEX NAME)



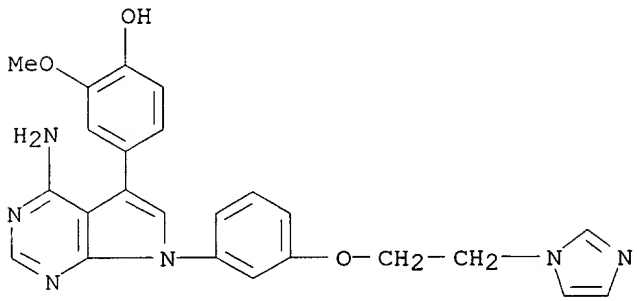
RN 178909-66-5 CAPLUS
CN Ethanol, 2-[3-[4-amino-5-(3-methoxyphenyl)-7H-pyrrolo[2,3-d]pyrimidin-7-yl]phenoxy]- (9CI) (CA INDEX NAME)



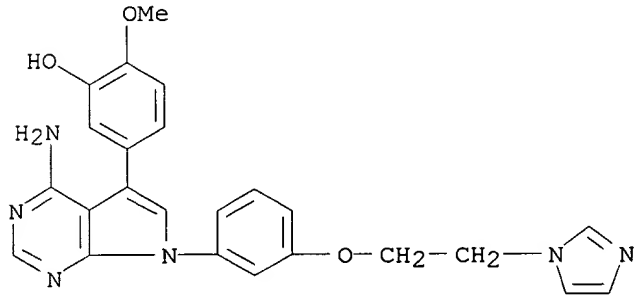
RN 178909-84-7 CAPLUS
CN Phenol, 4-(4-amino-7-phenyl-7H-pyrrolo[2,3-d]pyrimidin-5-yl)-2-methoxy- (9CI) (CA INDEX NAME)



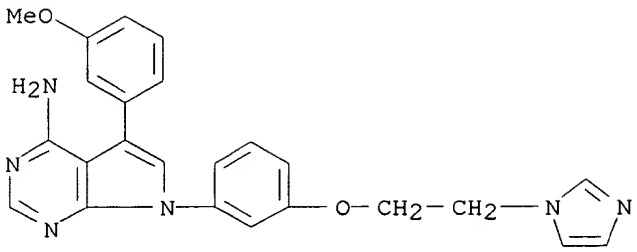
RN 178909-97-2 CAPLUS
CN Phenol,
4-[4-amino-7-[3-[2-(1H-imidazol-1-yl)ethoxy]phenyl]-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxy- (9CI) (CA INDEX NAME)



RN 178909-98-3 CAPLUS
CN Phenol,
5-[4-amino-7-[3-[2-(1H-imidazol-1-yl)ethoxy]phenyl]-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-methoxy- (9CI) (CA INDEX NAME)



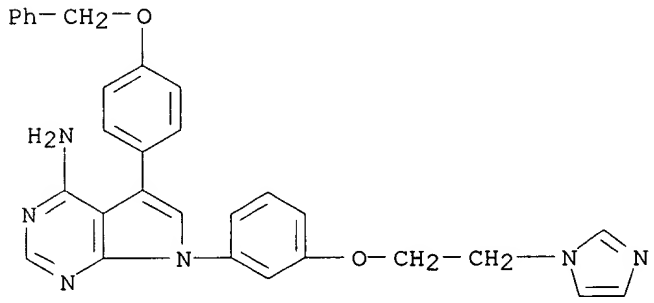
RN 178910-00-4 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 7-[3-[2-(1H-imidazol-1-yl)ethoxy]phenyl]-5-(3-methoxyphenyl)- (9CI) (CA INDEX NAME)



IT **178910-31-1P 178910-37-7P 178910-38-8P**
178910-39-9P 178910-40-2P 178910-41-3P
178910-42-4P 178910-43-5P 178910-44-6P
178910-45-7P 178910-46-8P 178910-47-9P
178910-48-0P 178910-49-1P 178910-50-4P
178910-51-5P 178910-52-6P 178910-53-7P
178910-54-8P 178910-55-9P 178910-56-0P
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation)
(prepn. of 4-aminopyrrolo[2,3-d]pyrimidines as inhibitors of the protein tyrosine kinase pp60c-src)

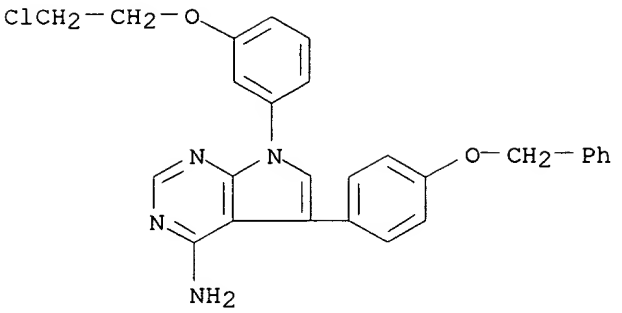
RN 178910-31-1 CAPLUS

CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 7-[3-[2-(1H-imidazol-1-yl)ethoxy]phenyl]-5-[4-(phenylmethoxy)phenyl]- (9CI) (CA INDEX NAME)

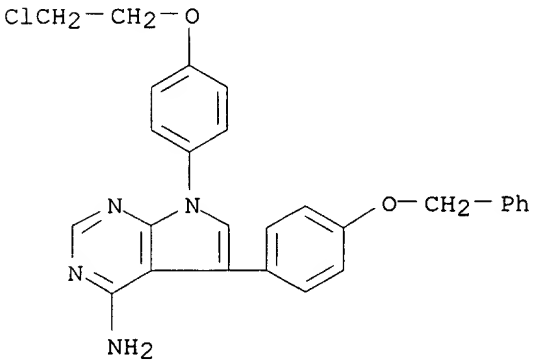


RN 178910-37-7 CAPLUS

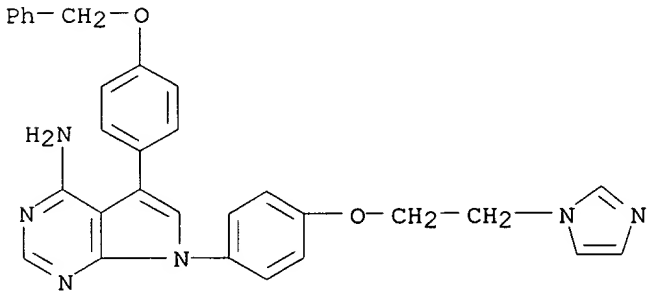
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 7-[3-(2-chloroethoxy)phenyl]-5-[4-(phenylmethoxy)phenyl]- (9CI) (CA INDEX NAME)



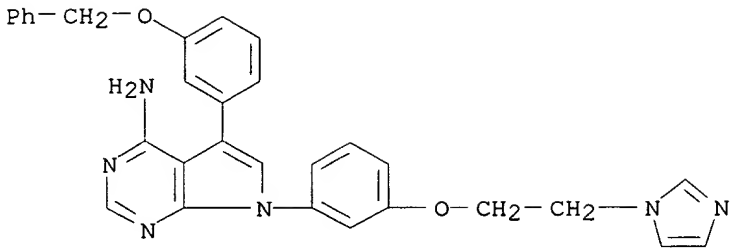
RN 178910-38-8 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 7-[4-(2-chloroethoxy)phenyl]-5-[4-(phenylmethoxy)phenyl]- (9CI) (CA INDEX NAME)



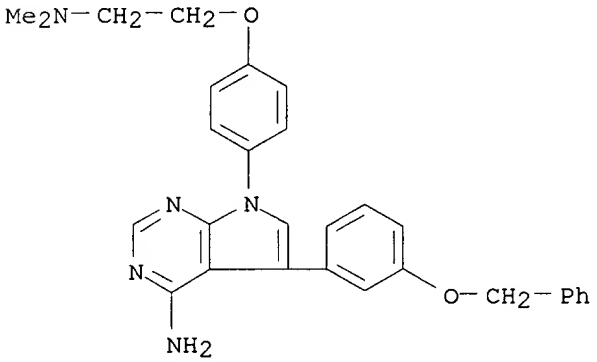
RN 178910-39-9 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 7-[4-[2-(1H-imidazol-1-yl)ethoxy]phenyl]-5-[4-(phenylmethoxy)phenyl]- (9CI) (CA INDEX NAME)



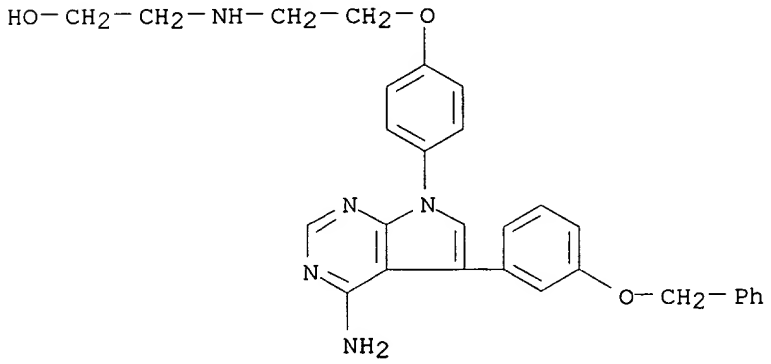
RN 178910-40-2 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 7-[3-[2-(1H-imidazol-1-yl)ethoxy]phenyl]-5-[3-(phenylmethoxy)phenyl]- (9CI) (CA INDEX NAME)



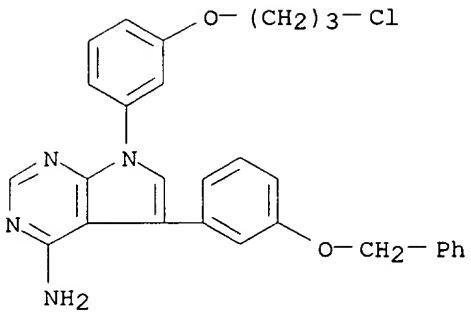
RN 178910-41-3 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 7-[4-[2-(dimethylamino)ethoxy]phenyl]-5-[3-(phenylmethoxy)phenyl]- (9CI) (CA INDEX NAME)



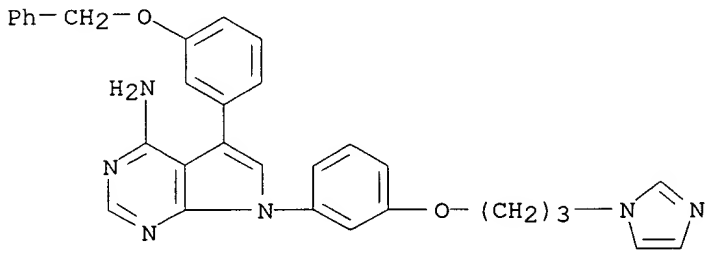
RN 178910-42-4 CAPLUS
CN Ethanol, 2-[[2-[4-[4-amino-5-[3-(phenylmethoxy)phenyl]-7H-pyrrolo[2,3-d]pyrimidin-7-yl]phenoxy]ethyl]amino]- (9CI) (CA INDEX NAME)



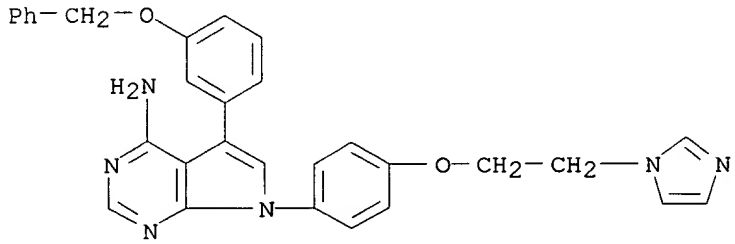
RN 178910-43-5 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 7-[3-(3-chloropropoxy)phenyl]-5-[3-(phenylmethoxy)phenyl]- (9CI) (CA INDEX NAME)



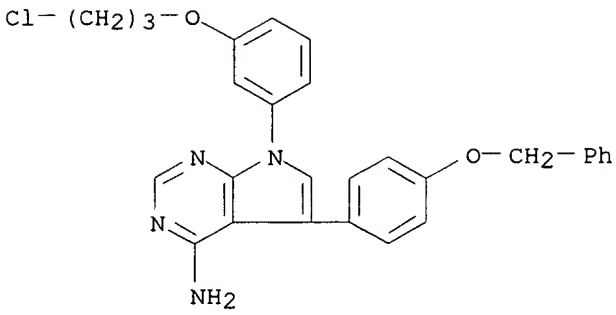
RN 178910-44-6 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 7-[3-[3-(1H-imidazol-1-yl)propoxy]phenyl]-5-[3-(phenylmethoxy)phenyl]- (9CI) (CA INDEX NAME)



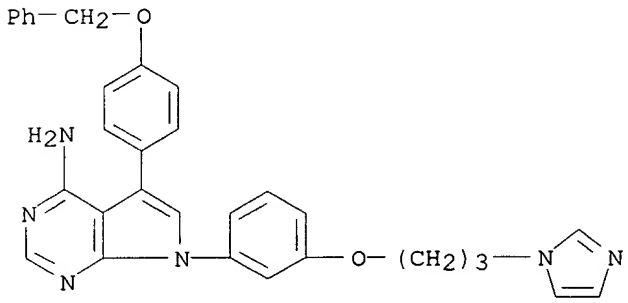
RN 178910-45-7 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 7-[4-[2-(1H-imidazol-1-yl)ethoxy]phenyl]-5-[3-(phenylmethoxy)phenyl]- (9CI) (CA INDEX NAME)



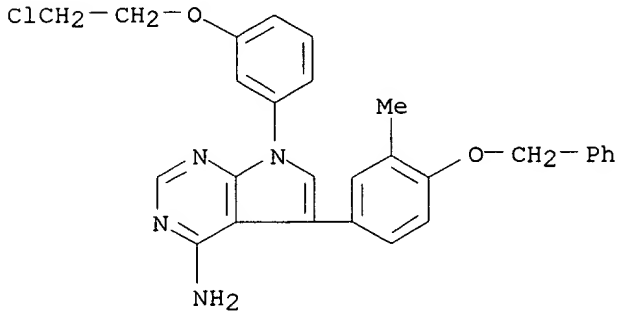
RN 178910-46-8 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 7-[3-(3-chloropropoxy)phenyl]-5-[4-(phenylmethoxy)phenyl]- (9CI) (CA INDEX NAME)



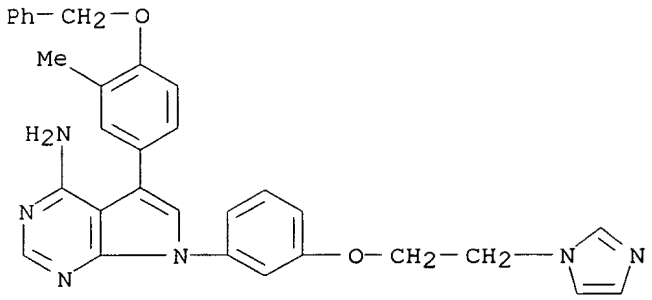
RN 178910-47-9 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 7-[3-[3-(1H-imidazol-1-yl)propoxy]phenyl]-5-[4-(phenylmethoxy)phenyl]- (9CI) (CA INDEX NAME)



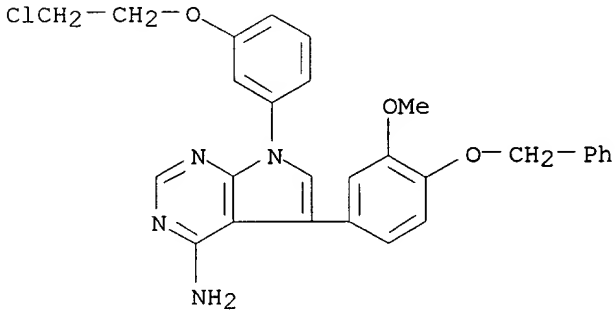
RN 178910-48-0 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 7-[3-(2-chloroethoxy)phenyl]-5-[3-methyl-4-(phenylmethoxy)phenyl]- (9CI) (CA INDEX NAME)



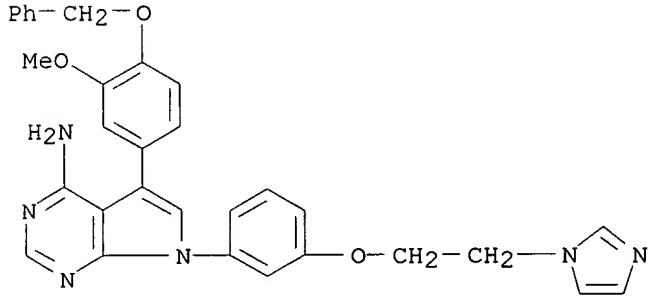
RN 178910-49-1 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 7-[3-[2-(1H-imidazol-1-yl)ethoxy]phenyl]-5-[3-methyl-4-(phenylmethoxy)phenyl]- (9CI) (CA INDEX NAME)



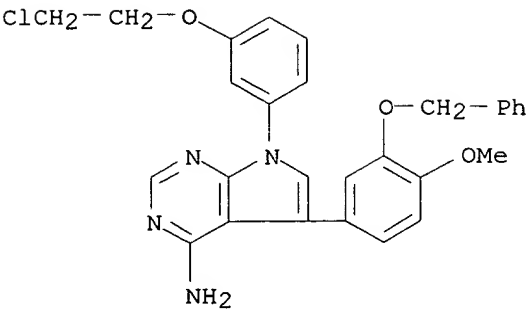
RN 178910-50-4 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 7-[3-(2-chloroethoxy)phenyl]-5-[3-methoxy-4-(phenylmethoxy)phenyl]- (9CI) (CA INDEX NAME)



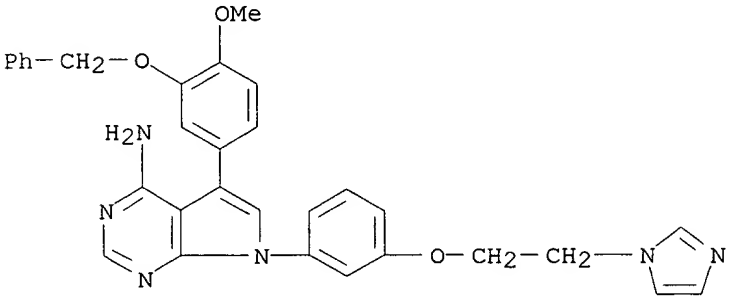
RN 178910-51-5 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 7-[3-[2-(1H-imidazol-1-yl)ethoxy]phenyl]-5-[3-methoxy-4-(phenylmethoxy)phenyl]- (9CI) (CA INDEX NAME)



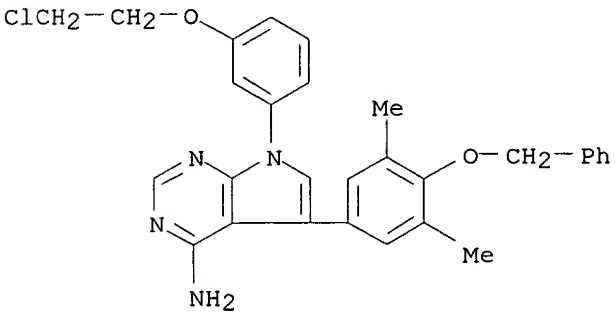
RN 178910-52-6 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 7-[3-(2-chloroethoxy)phenyl]-5-[4-methoxy-3-(phenylmethoxy)phenyl]- (9CI) (CA INDEX NAME)



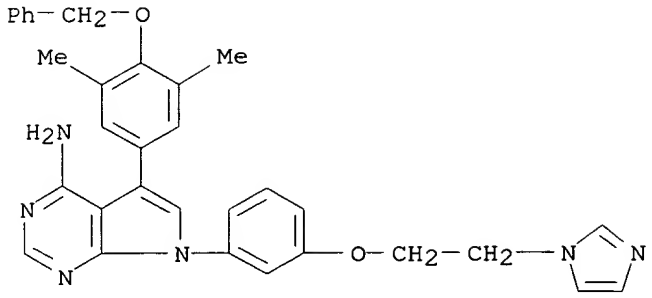
RN 178910-53-7 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 7-[3-[2-(1H-imidazol-1-yl)ethoxy]phenyl]-5-[4-methoxy-3-(phenylmethoxy)phenyl]- (9CI) (CA INDEX NAME)



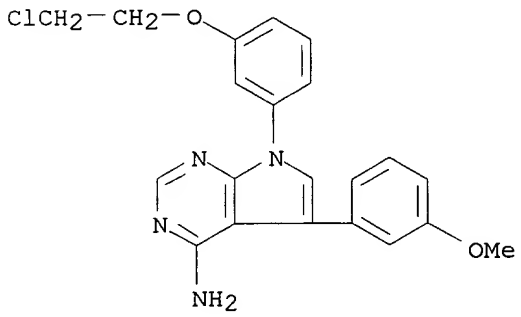
RN 178910-54-8 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 7-[3-(2-chloroethoxy)phenyl]-5-[3,5-dimethyl-4-(phenylmethoxy)phenyl]- (9CI) (CA INDEX NAME)



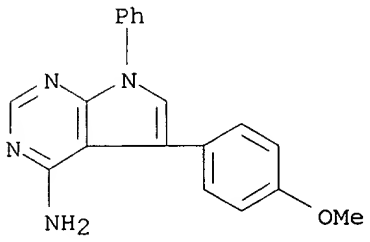
RN 178910-55-9 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 5-[3,5-dimethyl-4-(phenylmethoxy)phenyl]-7-[3-[2-(1H-imidazol-1-yl)ethoxy]phenyl]- (9CI) (CA INDEX NAME)



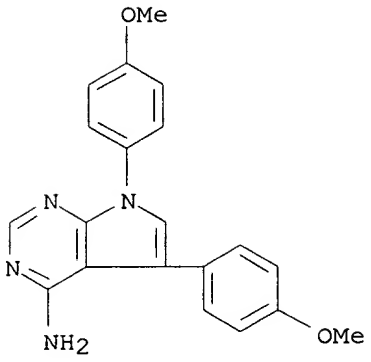
RN 178910-56-0 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 7-[3-(2-chloroethoxy)phenyl]-5-(3-methoxyphenyl)- (9CI) (CA INDEX NAME)



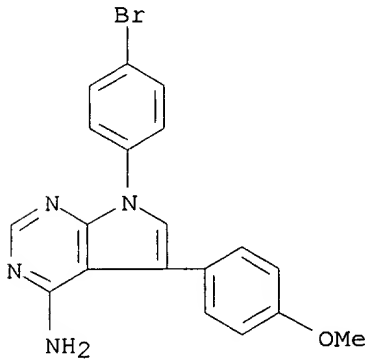
L6 ANSWER 10 OF 10 CAPLUS COPYRIGHT 2000 ACS
AN 1989:439290 CAPLUS
DN 111:39290
TI Synthesis and biological activity of pyrrolo[2,3-d]pyrimidines
AU Dave, Chaitanya G.; Shah, P. R.; Upadhyaya, S. P.; Gandhi, T. P.; Patel, R. B.
CS Dep. Chem., St. Xavier's Coll., Ahmedabad, 380 009, India
SO Indian J. Chem., Sect. B (1988), 27B(8), 778-80
CODEN: IJSBDB; ISSN: 0376-4699
DT Journal
LA English
OS CASREACT 111:39290
AB 2-Amno-3-pyrrolecarbonitriles were treated with HCONH2 to give aminopyrrolopyrimidines I [R1 = Ph, tolyl, anisyl, halophenyl; R2 = H, or R2R3 = (CH2)4; R3 = Ph, anisyl, ClC6H4, Me, tolyl]. Most I showed bactericidal, analgesic, antiinflammatory, antihistaminic, anticholinergic, anticonvulsant, and antihypertensive activity. Also prepd., from CS2, were pyrrolopyrimidines II.
IT **121405-30-9P**
RL: SPN (Synthetic preparation); PREP (Preparation)
(prepn. and anticholinergic activity of)
RN 121405-30-9 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 5-(4-methoxyphenyl)-7-phenyl- (9CI) (CA INDEX NAME)



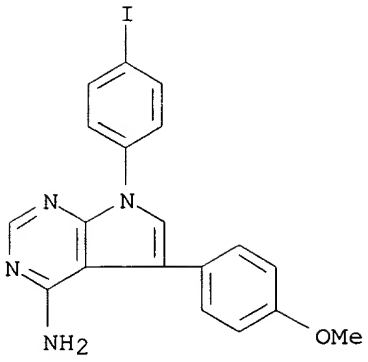
IT **121405-31-0P 121405-32-1P**
RL: BAC (Biological activity or effector, except adverse); SPN (Synthetic preparation); BIOL (Biological study); PREP (Preparation)
(prepn. and pharmacol. activity of)
RN 121405-31-0 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 5,7-bis(4-methoxyphenyl)- (9CI) (CA INDEX NAME)



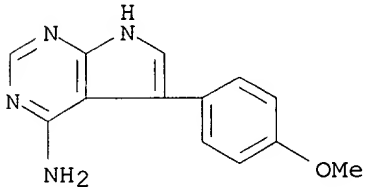
RN 121405-32-1 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine,
7-(4-bromophenyl)-5-(4-methoxyphenyl)-
(9CI) (CA INDEX NAME)



IT **121405-33-2P 121405-37-6P**
RL: SPN (Synthetic preparation); PREP (Preparation)
(prepn. of)
RN 121405-33-2 CAPLUS
CN 7H-Pyrrolo[2,3-d]pyrimidin-4-amine, 7-(4-iodophenyl)-5-(4-methoxyphenyl)-
(9CI) (CA INDEX NAME)

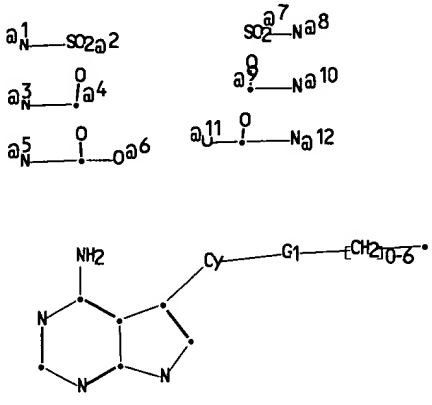


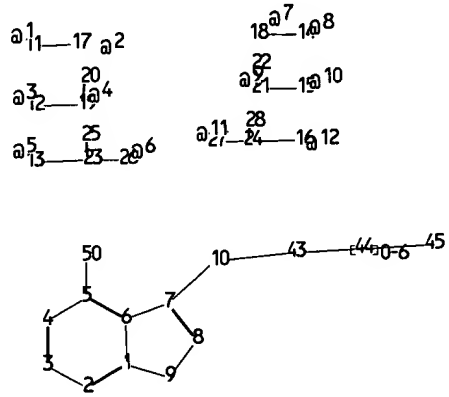
RN 121405-37-6 CAPLUS
CN 1H-Pyrrolo[2,3-d]pyrimidin-4-amine, 5-(4-methoxyphenyl)- (9CI) (CA INDEX NAME)



```
=> d his

     (FILE 'HOME' ENTERED AT 16:50:35 ON 06 MAR 2000)

     FILE 'REGISTRY' ENTERED AT 16:50:52 ON 06 MAR 2000
L1              STRUCTURE UPLOADED
L2           21 S L1 SSS SAM

     FILE 'STNGUIDE' ENTERED AT 16:52:10 ON 06 MAR 2000

     FILE 'REGISTRY' ENTERED AT 17:00:44 ON 06 MAR 2000
L3              STRUCTURE UPLOADED
L4            7 S L3 SSS SAM
L5          142 S L3 SSS FUL

     FILE 'CAPLUS' ENTERED AT 17:03:50 ON 06 MAR 2000
L6           10 S L5

     FILE 'CAOLD' ENTERED AT 17:05:44 ON 06 MAR 2000

=> s l5

L7            0 L5
```

C:\STNEXP4\QUERIES\09399083x.str





```
chain nodes :
    10  11  12  13  14  15  16  17  18  19  20  21  22  23  24  25  26  27  28  43  44  45  50
ring nodes :
    1  2  3  4  5  6  7  8  9
chain bonds :
    5-50  7-10  10-43  11-17  12-19  13-23  14-18  15-21  16-24  19-20  21-22  23-25  23-26
    24-27  24-28  43-44  44-45
ring bonds :
    1-2  1-6  1-9  2-3  3-4  4-5  5-6  6-7  7-8  8-9
exact/norm bonds :
    1-9  5-50  7-10  8-9  10-43  11-17  12-19  13-23  14-18  15-21  16-24  19-20  21-22  23-25
    23-26  24-27  24-28  43-44
exact bonds :
    6-7  7-8  44-45
normalized bonds :
    1-2  1-6  2-3  3-4  4-5  5-6
isolated ring systems :
    containing 1 :


G1:[*1-*2],[*3-*4],[*5-*6],[*7-*8],[*9-*10],[*11-*12]

Hydrogen count :
    3:>= minimum 1
Match level :
    1:Atom  2:Atom  3:Atom  4:Atom  5:Atom  6:Atom  7:Atom  8:Atom  9:Atom  10:Atom  11:CLASS
    12:CLASS  13:CLASS  14:CLASS  15:CLASS  16:CLASS  17:CLASS  18:CLASS  19:CLASS  20:CLASS
    21:CLASS  22:CLASS  23:CLASS  24:CLASS  25:CLASS  26:CLASS  27:CLASS  28:CLASS  43:CLASS
    44:CLASS  45:CLASS  50:CLASS
Generic attributes :
    10:
    Saturation              : Unsaturated
```